



UNITED STATES ENVIRONMENTAL PROTECTION AGENCY
REGION III
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SDMS DocID 2198336

DATE : October 13, 1992

SUBJECT : Region III Data QA Review

FROM : Cynthia E. Caporale *C Caporale*
Region III ESAT RPO (3ES31)

TO : Michael Taurino
Regional Project Manager (3HW73)

Attached is the organic data validation report for the Hoffman Landfill Site (Case 18347) completed by the Region III Environmental Services Assistance Team (ESAT) contractor under the direction of Region III ESD.

If you have any questions regarding this review, please call me.

Attachment

cc: Jennifer Woods, MD DOE
Edward Kantor, EMSL-LV
Regional CLP TPO: Tom Bennett

Region: IV Lab Code: COMPU

TID File: 03920418 Task 1514

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SITE ASSESSMENT
SECTION



Environmental Services Assistance Teams
Region 3

1419 Forest Drive, Suite 104
Annapolis, Maryland 21403

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DATE: October 1, 1992

SUBJECT: Organic Data Validation For Case 18347
Site: Hoffman Landfill

FROM: Hari Prasad
Organic Data Reviewer

Mahboobeh Mecanic
Senior Oversight Chemist

TO: Cynthia E. Caporale
ESAT Regional Project Officer

THROUGH: Dale S. Boshart
ESAT Team Manager

OVERVIEW

Case 18347 consisted of twelve (12) aqueous and eleven (11) soil samples submitted to Compuchem Laboratories, Inc. for volatile, semivolatile and pesticide/PCB analyses. The aqueous samples included one (1) field duplicate pair and one (1) field blank and the soil samples included one (1) field duplicate pair. The samples were analyzed as a Contract Laboratory Program (CLP) Routine Analytical Service (RAS).

SUMMARY

The samples were successfully analyzed for all target compounds except a few compounds in the semivolatile fraction. All other instrument and method sensitivities were according to the Contract Laboratory Program (CLP) Routine Analytical Service (RAS) protocol.

MAJOR PROBLEMS

- o In the semivolatile analyses, 3-nitroaniline had a relative response factor (RRF) less than 0.05 (< 0.05) in the initial calibration dated 6/15/92. The quantitation limits for this compound in the affected samples were qualified unusable ("R"). (See Table I in Appendix F.)
- o In the semivolatile analyses, sample CKY32 had the recovery for the acid surrogate, 2,4,6-tribromophenol, less than 10% and the quantitation limits for its acid compounds were qualified unusable ("R"). (See FORM-II SV-2 in Appendix F.)

MINOR PROBLEMS

- o The semivolatile extractions of samples CKY28, CKY31, CKY34, CKY35, CKY36 and CKY38 were performed thirteen (13) days and sample CKY29 seventeen (17) days from the date of sample collection. Although no technical holding time for the semivolatile extraction of soil samples has been established, the technical holding time of seven (7) days for the aqueous samples has been exceeded by six (6) and ten (10) days, respectively, in these samples. The aqueous sample holding time was applied and, therefore, the quantitation limits in these samples were qualified "UJ" and positive results were qualified "J", except when superseded by the "B" qualifier in the affected samples. (See Traffic Report in Appendix F.)
- o Several compounds failed precision criteria in the volatile and semivolatile initial and continuing calibrations. The positive results were qualified "J" except when superseded by the "B" qualifier. The quantitation limits were qualified "UJ" when the QC limits were grossly exceeded (%RSD or %D greater than 50%), except for the acid compound 4-nitrophenol in the semivolatile fraction of sample CKY32, where it was superseded by the "R" qualifier. (See Table I in Appendix F.)
- o During the pesticide/PCB analyses, positive results have been flagged "P" on Form I's when the %D between the two columns was greater than 25%. These results were qualified "J" on the data summary forms.

NOTES

- o During the semivolatile analysis, sample CKY18MSD failed both surrogate and spike recovery criteria due to an extraction error and no reanalysis was performed. No action was taken. (See Case Narrative.)
- o In the semivolatile analyses, aqueous samples CKY16, CKY19 and CKY20 had one (1) and sample CKY18 had two (2) (one acid and one base) surrogate recoveries above the QC limits and soil samples CKY32MS, CKY32MSD and CKY34 had one (1) surrogate recovery below the QC limit, but greater than 10%. No action was taken. (See FORM-II SV-1 and FORM-II SV-2 in Appendix F.)
- o In the pesticide/PCB analyses, aqueous samples CKY14 and CKY26 and soil samples CKY30, CKY32 and CKY36 had one (1) each of their surrogate recoveries below the QC limits but greater than 10%. No action was taken. (See FORM II-PEST in Appendix F.)
- o The maximum concentration of all compounds found in the analyses of the field and laboratory method blanks are listed below. Samples with concentrations of common

laboratory contaminants less than ten times (<10X) the blank concentration or with concentrations of other contaminants less than five times (<5X) the blank concentration, have been qualified "B" on the data summary forms.

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<u>COMPOUND</u>	<u>CONCENTRATION</u>
methylene chloride*	30 µg/Kg
acetone*	31 µg/Kg
phenol	3 J µg/L
diethylphthalate*	1 J µg/L
butylbenzylphthalate*	6 J µg/L
bis(2-ethylhexyl)phthalate*	10 µg/L
methoxychlor	0.59 J µg/Kg

* Common Laboratory Contaminant

- o Samples CKY16/CKY19 and CKY34/CKY38 were the field duplicate pairs analyzed in the aqueous and solid samples, respectively. Their results and precision estimates, excluding the blank contaminants, are listed in the table below:

<u>COMPOUND</u>	<u>CONCENTRATION (µg/L)</u>		<u>RPD</u>
	<u>CKY16</u>	<u>CKY19</u>	
chloroform	4 J	4 J	0
<u>CONCENTRATION (µg/Kg)</u>			
	<u>CKY34</u>	<u>CKY38</u>	
fluoranthene	39 J	61 J	44
pyrene	ND	55 J	IN
benzo(b)fluoranthene	ND	61 J	IN
benzo(a)pyrene	ND	61 J	IN
gamma-BHC	0.13 J	ND	IN
heptachlor	ND	0.094 J	IN
aldrin	0.58 J	0.37 J	44
heptachlor epoxide	0.30 J	0.26 J	14
endosulfan I	0.40 J	0.34 J	16
dieldrin	ND	0.33 J	IN
endrin	0.59 J	0.56 J	5
endosulfan II	0.31 J	ND	IN
4,4'-DDD	ND	1.3 J	IN
endrin ketone	0.15 J	ND	IN
alpha-chlordane	0.55 J	0.48 J	14
gamma-chlordane	0.67 J	0.51 J	27

RPD = Relative Percent Difference

ND = Not Detected

IN = Indeterminate

- Non-spiked compounds other than blank contaminants were detected in the pesticide/PCB analyses of samples CKY18, CKY32 and their matrix spike/matrix spike duplicate recoveries. Their results and precision estimates are as follows:

<u>COMPOUND</u>	<u>CONCENTRATION ($\mu\text{g/L}$)</u>			<u>%RSD</u>
	<u>CKY18</u>	<u>CKY18MS</u>	<u>CKY18MSD</u>	
endrin ketone	ND	0.0099 J	0.0074 J	24
<u>CONCENTRATION ($\mu\text{g/Kg}$)</u>				
	<u>CKY32</u>	<u>CKY32MS</u>	<u>CKY32MSD</u>	
benzo(b)fluoranthene	ND	ND	48 J	IN
benzo(k)fluoranthene	ND	ND	48 J	IN
delta-BHC	ND	0.093 J	0.22 J	81+
heptachlor epoxide	ND	0.21 J	ND	IN
endrin ketone	0.66 J	ND	1.2 J	58+
benzo(b)fluoranthene	ND	ND	48 J	IN
benzo(k)fluoranthene	ND	ND	48 J	IN

%RSD = Percent Relative Standard Deviation

+ = RPD instead of %RSD

ND = Not Detected

IN = indeterminate

- In the semivolatile and pesticide/PCB analyses of soil samples, GPC cleanup was performed. The dilution factor of two (2) required by this procedure was accounted for in the analytical procedures used by the laboratory and, therefore, is not reflected in the data summary forms.
- Sample weights other than thirty (30) grams were used in the semivolatile and pesticide/PCB analyses of several soil samples. The dilution factors on the data summary forms have been changed to reflect this variance, when significant.
- During the semivolatile analyses of samples CKY29 and CKY38 benzo(b/k)fluoranthene isomers coeluted and their results have been flagged "X" on Form I's. (See case narrative in Appendix F.)
- Tentatively Identified Compounds (TIC's) in Appendix D were reviewed and corrected during data validation. Several early eluting TICs were found during the semivolatile analyses due to the use of contaminated methylene chloride. Compounds identified as solvents, laboratory artifacts or blank contaminants were crossed off the TIC Form I's. (See case narrative in Appendix F.)

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All data for case 18347 were reviewed in accordance with the Functional Guidelines for Evaluating Organic Analyses with Modifications for use within Region III. The text of this report addresses only those problems affecting usability.

Attachments

- 1) Appendix A - Glossary of Data Qualifiers
- 2) Appendix B - Data Summary. These include:
 - (a) All positive results for target compounds with qualifier codes where applicable.
 - (b) All unusable detection limits (qualified "R").
- 3) Appendix C - Results as reported by the Laboratory for all target compounds.
- 4) Appendix D - Reviewed and corrected Tentatively Identified Compounds.
- 5) Appendix E - Organic Regional Data Assessment Summary.
- 6) Appendix F - Support Documentation.

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Appendix A
Glossary of Data Qualifiers

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GLOSSARY OF DATA QUALIFIER CODES (ORGANIC)

CODES RELATING TO IDENTIFICATION

(confidence concerning presence or absence of compounds)

U = Not detected. The associated number indicates approximate sample concentration necessary to be detected.

NO CODE = Confirmed identification.

B = Not detected substantially above the level reported in laboratory or field blanks.

R = Unreliable result. Analyte may or may not be present in the sample. Supporting data necessary to confirm result.

N = Tentative identification. Consider present. Special methods may be needed to confirm its presence or absence in future sampling efforts.

CODES RELATED TO QUANTITATION

(can be used for both positive results and sample quantitation limits):

J = Analyte present. Reported value may not be accurate or precise.

K = Analyte present. Reported value may be biased high. Actual value is expected to be lower.

L = Analyte present. Reported value may be biased low. Actual value is expected to be higher.

UJ = Not detected, quantitation limit may be inaccurate or imprecise.

UL = Not detected, quantitation limit is probably higher.

OTHER CODES

Q = No analytical result.

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Appendix B

Data Summary Forms

DATA SUMMARY FORM: VOLATILES I

Site Name: HOFFMAN LANDFILL

Case #: 18347 Sampling Date(s): 6/23/92

WATER SAMPLES

(μ g/L)

To calculate sample quantitation limit:
 • Dilution Factor)

CROI as contract Required Quantitation Limit

DATA SUMMARY FORM: VOLATILES

Site Name: HOFFMAN LANDFILL

Case #: 18347 Sampling Date(s): 6/23/92

WATER SAMPLES

(µg/L)

To calculate sample quantitation limit:
(CROL * Dilution Factor)

	Sample No.	Dilution Factor	Location	CKY14	CKY15	CKY16	CKY17	CKY18	CKY19	CKY20	CKY22	CKY23
CROL	COMPOUND											
10	*1,2-Dichloropropane											
10	Cis-1,3-Dichloropropane											
10	Trichloroethene											
10	Dibromochloromethane											
10	1,1,2-Trichlorethane											
10	*Benzene											
10	Trans-1,3-Dichloropropene											
10	Bromoform											
10	4-Methyl-2-pentanone				VJ		VJ					
10	2-Hexanone											
10	*Tetrachloroethene											
10	1,1,2,2-Tetrachloroethane											
10	*Toluene											
10	*Chlorobenzene											
10	*Ethylbenzene											
10	*Styrene											
10	*Total Xylenes											

CROL = Contract Required Quantitation Limit

*Action Level Exists

SEE NARRATIVE FOR CODE DEFINITIONS

revised 07/90

DATA SUMMARY FORM: VOLATILES 1

Site Name: HOFFMAN LANDFILL

WATER SAMPLES

(μ g/L)

Case #: 18347 Sampling Date(s): 6/23/92

To calculate sample quantitation (CRQL * Dilution limit: factor)

CROL = Contract Required Quantitation Limit

DATA SUMMARY FORM: VOLATILES 2

Site Name: HOFFMAN LANDFILL

WATER SAMPLES
(kg/l)

Case #: 18347 Sampling Date(s): 6/23/92

To calculate	sample (C <small>oncentration</small>)	quantitation • dilution	limit: factor)
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CROL = Contract Required Quantitation Limit

***Action Level Exists**

SEE NARRATIVE FOR CODE DEFINITIONS

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DATA SUMMARY FORM: VOLATILES 1

Site Name: HOFFMAN LANDFILL

SOIL SAMPLES
($\mu\text{g}/\text{Kg}$)

Case #: 18347 Sampling Date(s): 6/23/92

To calculate sample quantitation limit:
 (CQL * dilution factor) / ((100 - % moisture)/100)

Sample No.
Dilution Factor
% Moisture
Location

	CKY28	CKY29	CKY30	CKY31	CKY32	CKY33	CKY34	CKY35	CKY36
	1.0	1.0	1.0	1.0	1.0	1.0	1.0	1.0	1.0
	19	17	40	23	18	11	13	3	10
	SFD1	SFD2	SFD3	SFD4	S1	S2			

CQL COMPOUND

10	Chloromethane								
10	Bromomethane								
10	Vinyl Chloride								
10	Chloroethane								
10	Methylene Chloride	30 B	78 B	84 B	100 B	33 B	31 B	48 B	87 B
10	Acetone	20 B	32 B	42 B	110 B	23 B	23 B	27 B	27 B
10	Carbon Disulfide								
10	1,1-Dichloroethene								
10	1,1-Dichloroethane								
10	Total 1,2-Dichloroethene								
10	Chloroform								
10	1,2-Dichloroethane								
10	2-Butanone								
10	1,1,1-Trichloroethane								
10	Carbon Tetrachloride								
10	Bromodichloromethane								

Field Dup.
of CKY38

DATA SUMMARY FORM: VOLATILES 2

Site Name: HOFFMAN LANDFILL

SOIL SAMPLES

(µg/Kg)

Case #: 18347 Sampling Date(s): 6/23/92

To calculate sample quantitation limit:
 $(CRQL * Dilution Factor) / ((100 - X moisture)/100)$

CRQL	COMPOUND	CKY28	CKY29	CKY30	CKY31	CKY32	CKY33	CKY34	CKY35	CKY36
		Dilution Factor	X Moisture	Location						
10	1,2-Dichloropropane									
10	Cis-1,3-Dichloropropene									
10	Trichloroethene									
10	Dibromochloromethane									
10	1,1,2-Trichloroethane									
10	Benzene									
10	Trans-1,3-Dichloropropene									
10	Bromoform									
10	4-Methyl-2-pentanone									
10	2-Hexanone									
10	Tetrachloroethene									
10	1,1,2,2-Tetrachloroethane									
10	Toluene	1 J								
10	Chlorobenzene									
10	Ethylbenzene									
10	Styrene									
10	Total Xylenes									

DATA SUMMARY FORM: VOLATILES 1

Site Name: HOFFMAN LANDFILLSOIL SAMPLES
($\mu\text{g}/\text{Kg}$)

Case #: 18347 Sampling Date(s): 6/23/92

To calculate sample quantitation limit:
(CRQL * Dilution factor) / ((100 - % moisture)/100)

Sample No.	CKY37	CKY38
Dilution Factor	1.0	1.0
% Moisture	8	12
Location	Field Dup. of CKY34	

CRQL	COMPOUND
10	Chloromethane
10	Bromomethane
10	Vinyl Chloride
10	Chloroethane
10	Methylene Chloride
10	Acetone
10	Carbon Disulfide
10	1,1-Dichloroethene
10	1,1-Dichloroethane
10	Total 1,2-Dichloroethene
10	Chloroform
10	1,2-Dichloroethane
10	2-Butanone
10	1,1,1-Trichloroethane
10	Carbon Tetrachloride
10	Bromodichloromethane

CRQL = Contract Required Quantitation Limit

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Revised 07/90

DATA SUMMARY FORM: VOLATILES 2

Site Name: HOFFMAN LANDFILL

SOIL SAMPLES

Case #: 18347 Sampling Date(s): 6/23/92

To calculate sample quantitation limits
 (CQL) = Dilution Factor / ((100 - x moisture)/100)

CROL = Contract Required Quantitation Limit

SEE NARRATIVE FOR CODE DEFINITIONS

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DATA SUMMARY FORM: B N A S 1

Site Name: HOFFMANN LANDFILL

Case #: 18347 Sampling Date(s): 6/23/92

WATER SAMPLES
($\mu\text{g/L}$)To calculate sample quantitation
(CROL * Dilution)

(limit factor)

CROL	COMPOUND	CKY14	CKY15	CKY16	CKY17	CKY18	CKY19	CKY20	CKY22	CKY23
		1.0 BLK1 Field Blank	1.0 GW1	1.0 GW2	1.0 GW3 Field Dup. of CKY19	1.0 GW4	1.0 GW5 Field Dup. of CKY16	1.0 GW6	1.0 SW1	1.0 SW2
10	Phenol									
10	bis(2-Chloroethyl)ether									
10	2-Chlorophenol									
10	*1,3-Dichlorobenzene									
10	*1,4-Dichlorobenzene									
10	1,2-Dichlorobenzene									
10	2-Methylphenol									
10	2,2'-Oxybis(1-chloropropane)									
10	4-Methylphenol									
10	N-Nitroso-di-n-propylamine									
10	Hexachloroethane									
10	Nitrobenzene									
10	Isophorone									
10	2-Nitrophenol									
10	2,4-Dimethylphenol									
10	bis(2-Chloroethoxy)methane									
10	2,4-Dichlorophenol									
10	1,2,4-Trichlorobenzene									
10	Naphthalene									
10	4-Chloroaniline		VJ	VJ	VJ	VJ	VJ	VJ	VJ	VJ

2

B

DATA SUMMARY FORM: B N A S 2

Site Name: HOFFMAN LANDFILL

Site #: 18347 Sampling Date(s): 6/23/92

WATER SAMPLES
($\mu\text{g/L}$)To calculate sample quantitation limits
(CRQL * Dilution Factor)

Sample No.	CKY14	CKY15	CKY16	CKY17	CKY18	CKY19	CKY20	CKY22	CKY25
Dilution Factor	1.0	1.0	1.0	1.0	1.0	1.0	1.0	1.0	1.0
Location	BLK1	GW1	GW2	GW3	GW4	GW5	GW6	SW1	SW2
Field Blank				Field Dup. of CKY19			Field Dup. of CKY16		
CRQL	COMPOUND								
10	Hexachlorobutadiene								
10	4-Chloro-3-methylphenol								
10	2-Methylnaphthalene								
10	Hexachlorocyclopentadiene								
10	2,4,6-Trichlorophenol								
25	2,4,5-Trichlorophenol								
10	2-Chloronaphthalene								
25	2-Nitroaniline								
10	Dimethylphthalate								
10	Acenaphthylene								
10	2,6-Dinitrotoluene								
25	3-Nitroaniline								
10	Acenaphthene								
25	2,4-Dinitrophenol								
25	4-Nitrophenol								
10	Dibenzofuran								
10	2,4-Dinitrotoluene								
10	Diethylphthalate								
10	4-Chlorophenyl-phenylether								
10	Fluorene								
25	4-Nitroaniline								
25	4,6-Dinitro-2-methylphenol								

RQL = Contract Required Quantitation Limit

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DATA SUMMARY FORM: B N A S 3

Site Name: HOFFMAN LANDFILL

Case #: 18347 Sampling Date(s): 6/23/92

WATER SAMPLES
($\mu\text{g/L}$)

To calculate sample quantitation (CRQL * Dilution Factor)

CRQL	COMPOUND	CKY14	CKY15	CKY16	CKY17	CKY18	CKY19	CKY20	CKY22	CKY23
		Dilution Factor	Sample No.	Location	1.0	1.0	1.0	1.0	1.0	1.0
10	N-Nitrosodiphenylamine									
10	4-Bromophenyl-phenylether									
10	*Hexachlorobenzene									
25	*Pentachlorophenol									
10	Phenanthrene									
10	Anthracene									
10	Carbazole									
10	Di-n-butylphthalate									
10	Fluoranthene									
10	Pyrene									
10	Butylbenzylphthalate									
10	3,3'-Dichlorobenzidine									
10	Benzo(a)anthracene									
10	Chrysene									
10	bis(2-Ethylhexyl)phthalate	2 B	2 B	1 B	3 B	3 B	2 B	6 B	6 B	3 B
10	Di-n-octylphthalate									
10	Benzo(b)fluoranthene									
10	Benzo(k)fluroanthene									
10	Benzo(a)pyrene									
10	Indeno(1,2,3-cd)pyrene									
10	Dibenz(a,h)anthracene									
10	Benzo(g,h,i)perylene									

DATA SUMMARY FORM: B N A S 1

Site Name: HOFFMAN LANDFILLCase #: 18347 Sampling Date(s): 6/23/92WATER SAMPLES
($\mu\text{g/L}$)To calculate sample quantitation limit:
(CRQL * Dilution Factor)

CRQL	COMPOUND	CKY24	CKY25	CKY26								
		Sample No.	Dilution Factor	Location								
10	Phenol											
10	bis(2-Chloroethyl)ether											
10	2-Chlorophenol											
10	*1,3-Dichlorobenzene											
10	*1,4-Dichlorobenzene											
10	1,2-Dichlorobenzene											
10	2-Methylphenol											
10	2,2'-Oxybis(1-chloropropane)											
10	4-Methylphenol											
10	N-Nitroso-di-n-propylamine											
10	Hexachloroethane											
10	Nitrobenzene											
10	Isophorone											
10	2-Nitrophenol											
10	2,4-Dimethylphenol											
10	bis(2-Chloroethoxy)methane											
10	2,4-Dichlorophenol											
10	1,2,4-Trichlorobenzene											
10	Naphthalene											
10	4-Chloroaniline											

CRQL = Contract Required Quantitation Limit

DATA SUMMARY FORM: B N A S 2

Site Name: HOFFMAN LANDFILL

Case #: 18347 Sampling Date(s): 5/23/92

WATER SAMPLES

(µg/L)

To calculate sample quantitation limits
(CQL * Dilution Factor)

CQL	COMPOUND	CKY24	CKY25	CKY26								
		Dilution Factor	Location									
10	Hexachlorobutadiene											
10	4-Chloro-3-methylphenol											
10	2-Methylnaphthalene											
10	Hexachlorocyclopentadiene											
10	2,4,6-Trichlorophenol											
25	2,4,5-Trichlorophenol											
10	2-Chloronaphthalene											
25	2-Nitroaniline											
10	Dimethylphthalate											
10	Acenaphthylene											
10	2,6-Dinitrotoluene											
25	3-Nitroaniline	R	R	R								
10	Acenaphthene											
25	2,4-Dinitrophenol											
25	4-Nitrophenol											
10	Dibenzofuran											
10	2,4-Dinitrotoluene											
10	Diethylphthalate											
10	4-Chlorophenyl-phenylether											
10	Fluorene											
25	4-Nitroaniline											
25	4,6-Dinitro-2-methylphenol											

CQL = Contract Required Quantitation Limit

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DATA SUMMARY FORM: B N A S 3

WATER SAMPLES

(µg/L)

Site Name: HOFFMAN LANDFILL
 Case #: 18347 Sampling Date: 6/23/92

To calculate sample quantitation
 (CRQL * Dilution Factor)

Sample No. Dilution Factor Location	CKY24	CKY25	CKY26												
	1.0 SW3	1.0 SW4	1.0 LTL												
CRQL COMPOUND															
10	N-Nitrosodiphenylamine														
10	4-Bromophenyl-phenylether														
10	*Hexachlorobenzene														
25	*Pentachlorophenol														
10	Phenanthrene														
10	Anthracene														
10	Carbazole														
10	Di-n-butylphthalate														
10	Fluoranthene														
10	Pyrene														
10	Butylbenzylphthalate														
10	3,3'-Dichlorobenzidine	UT	UT	UT											
10	Benzo(a)anthracene														
10	Chrysene														
10	bis(2-Ethylhexyl)phthalate	3 B	1 B	2 B											
10	Di-n-octylphthalate														
10	Benzo(b)fluoranthene														
10	Benzo(k)fluroanthene														
10	Benzo(a)pyrene														
10	Indeno(1,2,3-cd)pyrene														
10	Dibenz(a,h)anthracene														
10	Benzo(g,h,i)perylene														

DATA SUMMARY FORM: B N A S 1

SOIL SAMPLES
($\mu\text{g}/\text{Kg}$)

Site Name: HOFFMAN LANDFILL

Case #: 18347 Sampling Date: 6/23/92

To calculate sample quantitation limit:

(CRQL * Dilution Factor) / ((1 - % moisture)/100)

CROL	COMPOUND	CKY28	CKY29	CKY30	CKY31	CKY32	CKY33	CKY34	CKY35	CKY36
		1.0	1.0	1.0	1.0	1.0	1.0	1.0	1.0	1.0
		% Moisture	19	17	40	23	18	11	13	3
330	Phenol		UJ			UJ		UJ	UJ	UJ
330	bis(2-Chloroethyl)ether									
330	2-Chlorophenol									
330	1,3-Dichlorobenzene									
330	1,4-Dichlorobenzene									
330	1,2-Dichlorobenzene									
330	2-Methylphenol									
330	2,2'-Oxybis(1-chloropropane)					UJ		UJ	UJ	UJ
330	4-Methylphenol					62 J		UJ	UJ	
330	N-Nitroso-di-n-propylamine									
330	Hexachloroethane									
330	Nitrobenzene									
330	Isophorone									
330	2-Nitrophenol									
330	2,4-Dimethylphenol									
330	bis(2-Chloroethoxy)methane									
330	2,4-Dichlorophenol									
330	1,2,4-Trichlorobenzene									
330	Naphthalene		V							
330	4-Chloroaniline					↓			↓	V

CROL = Contract Required Quantitation Limit

SEE NARRATIVE FOR CODE DEFINITIONS
revised 07/90

DATA SUMMARY FORM: B N A S 2

Site Name: HOFFMAN LANDFILL

Case #: 18347 Sampling Date(s): 6/23/92

SOIL SAMPLES
($\mu\text{g}/\text{Kg}$)

To calculate sample quantitation limits:
 $(\text{CQL} \times \text{Dilution Factor}) / ((100 - \% \text{ moisture})/100)$

CQL	COMPOUND	CKY28	CKY29	CKY30	CKY31	CKY32	CKY33	CKY34	CKY35	CKY36
		1.0	1.0	1.0	1.0	1.0	1.0	1.0	1.0	1.0
330	Hexachlorobutadiene	UJ				UJ		UJ	UJ	UJ
330	4-Chloro-3-methylphenol									
330	2-Methylnaphthalene		92	JT						
330	Hexachlorocyclopentadiene									
330	2,4,6-Trichlorophenol									
800	2,4,5-Trichlorophenol									
330	2-Chloronaphthalene									
800	2-Nitroaniline									
330	Dimethylphthalate									
330	Acenaphthylene									
330	2,6-Dinitrotoluene									
800	3-Nitroaniline									
330	Acenaphthene									
800	2,4-Dinitrophenol									
800	4-Nitrophenol									
330	Dibenzofuran									
330	2,4-Dinitrotoluene									
330	Diethylphthalate									
330	4-Chlorophenyl-phenylether									
330	Fluorene									
800	4-Nitroaniline			UJ		VJ		VJ		UJ
800	4,6-Dinitro-2methylphenol	V			V			V	V	V

CQL = Contract Required Quantitation Limit

SEE NARRATIVE FOR CODE DEFINITIONS

revised 07/90

DATA SUMMARY FORM: B N A S 3

Site Name: HOFFMAN LANDFILL

Case #: 18347 Sampling Date(s): 6/23/92

SOIL SAMPLES
($\mu\text{g}/\text{Kg}$)

To calculate sample quantitation limit:
 $(\text{CRQL} * \text{Dilution Factor}) / ((100 - \% \text{ moisture})/100)$

Sample No.	Dilution Factor	CKY28	CKY29	CKY30	CKY31	CKY32	CKY33	CKY34	CKY35	CKY36
		1.0	1.0	1.0	1.0	1.0	1.0	1.0	1.0	1.0
X Moisture	19	17	40	23	18	11	13	3	10	10
Location	SFD1	SFD2	SFD3	SFD4	S1	S2	S3	S4	S5	S6
CRQL	COMPOUND									
330	N-Nitrosodiphenylamine		VJ			VJ		VJ	VJ	VJ
330	4-Bromophenyl-phenylether									
330	Hexachlorobenzene									
800	Pentachlorophenol									
330	Phenanthrene	130	J	56 J						
330	Anthracene									
330	Carbazole									
330	Di-n-butylphthalate									
330	Fluoranthene	120	J					39 J		
330	Pyrene	85	J					VJ		
330	Butylbenzylphthalate	46	B	240 B		180 B	160 B			
330	3,3'-Dichlorobenzidine									
330	Benzo(a)anthracene	110	J							
330	Chrysene	120	J							
330	bis(2-Ethylhexyl)phthalate	93	B				38 B			
330	Di-n-octylphthalate									
330	Benzo(b)fluoranthene	250	J							
330	Benzo(k)fluoranthene	250	J							
330	Benzo(a)pyrene	120	J							
330	Indeno(1,2,3-cd)pyrene	64	J							
330	Dibenz(a,h)anthracene									
330	Benzo(g,h)perylene	V				V		V	V	V

CRQL = Contract Required Quantitation Limit

SEE NARRATIVE FOR CODE DEFINITIONS

revised 07/90

DATA SUMMARY FORM: B N A S 1

Site Name: HOFFMAN LANDFILL

Case #: 18347 Sampling Date: 6/23/92

SOIL SAMPLES
($\mu\text{g}/\text{Kg}$)

To calculate sample quantitation limit:

(CQL * Dilution Factor) / ((1 - % moisture)/100)

CQL	COMPOUND	Sample No.													
		Dilution Factor	% Moisture												
330	Phenol	CK437	CK438												
330	bis(2-Chloroethyl)ether	1.0	1.0												
330	2-Chlorophenol	8	12												
330	1,3-Dichlorobenzene	S6	S7												
		Field Dup. of CK434													
330	1,4-Dichlorobenzene														
330	1,2-Dichlorobenzene														
330	2-Methylphenol														
330	2,2'-Oxybis(1-chloropropane)														
330	4-Methylphenol														
330	N-Nitroso-di-n-propylamine														
330	Hexachloroethane														
330	Nitrobenzene														
330	Isophorone														
330	2-Nitrophenol														
330	2,4-Dimethylphenol														
330	bis(2-Chloroethoxy)methane														
330	2,4-Dichlorophenol														
330	1,2,4-Trichlorobenzene														
330	Naphthalene														
330	4-Chloroaniline														

CQL = Contract Required Quantitation Limit

SEE NARRATIVE FOR CODE DEFINITIONS

revised 07/90

DATA SUMMARY FORM: B N A S 2

Site Name: HOFFMAN LANDFILL
 Case #: 18347 Sampling Date(s): 6/23/92

SOIL SAMPLES
($\mu\text{g}/\text{Kg}$)

To calculate sample quantitation limit:
 $(\text{CQL} * \text{Dilution Factor}) / ((100 + \% \text{ moisture})/100)$

CQL	COMPOUND	Sample No.	CKY37	CKY38	1.0	1.0	8	12	S6	S7												
330	Hexachlorobutadiene																					
330	4-Chloro-3-methylphenol																					
330	2-Methylnaphthalene																					
330	Hexachlorocyclopentadiene																					
330	2,4,6-Trichlorophenol																					
800	2,4,5-Trichlorophenol																					
330	2-Chloronaphthalene																					
800	2-Nitroaniline																					
330	Dimethylphthalate																					
330	Acenaphthylene																					
330	2,6-Dinitrotoluene																					
800	3-Nitroaniline																					
330	Acenaphthene																					
800	2,4-Dinitrophenol																					
800	4-Nitrophenol																					
330	Dibenzofuran																					
330	2,4-Dinitrotoluene																					
330	Diethylphthalate																					
330	4-Chlorophenyl-phenylether																					
330	Fluorene																					
800	4-Nitroaniline			VJ																		
800	4,6-Dinitro-2methylphenol																					

CQL = Contract Required Quantitation Limit

SEE NARRATIVE FOR CODE DEFINITIONS

revised 07/90

DATA SUMMARY FORM: B N A S 3

Site Name: HOFFMAN LANDFILL
 Case #: 18347 Sampling Date(s): 6/23/92

SOIL SAMPLES
($\mu\text{g/Kg}$)

To calculate sample quantitation limit:

$$(\text{CRQL} * \text{Dilution Factor}) / ((100 - \% \text{ moisture})/100)$$

CRQL	COMPOUND	Sample No.	Dilution Factor	% Moisture	Location	SOIL SAMPLES ($\mu\text{g/Kg}$)														
						CKY37	CKY38	1.0	1.0	8	12	S6	S7	Field Dip	cf CKY34	1.0	1.0	1.0	1.0	1.0
330	N-Nitrosodiphenylamine																			
330	4-Bromophenyl-phenylether																			
330	Hexachlorobenzene																			
800	Pentachlorophenol																			
330	Phenanthrene																			
330	Anthracene																			
330	Carbazole																			
330	Di-n-butylphthalate																			
330	Fluoranthene																			
330	Pyrene																			
330	Butylbenzylphthalate	130	B																	
330	3,3'-Dichlorobenzidine																			
330	Benzo(a)anthracene																			
330	Chrysene																			
330	bis(2-Ethylhexyl)phthalate																			
330	Di-n-octylphthalate																			
330	Benzo(b)fluoranthene																			
330	Benzo(k)fluoranthene																			
330	Benzo(a)pyrene																			
330	Indeno(1,2,3-cd)pyrrene																			
330	Dibenz(a,h)anthracene																			
330	Benzo(g,h)perylene																			

CRQL = Contract Required Quantitation Limit

SEE NARRATIVE FOR CODE DEFINITIONS
 revised 07/90

DATA SUMMARY FORM: PESTICIDES AND PCB'S

Site Name: HOFFMAN LANDFILLCase #: 18347 Sampling Date(s): 6/23/92WATER SAMPLES
($\mu\text{g/l}$)To calculate sample quantitation limits:
(CRQL * Dilution Factor)

Sample No.	CKY14	CKY15	CKY 16	CKY 17	CKY 18	CKY 19	CKY 20	CKY22	CKY23
Dilution Factor	1.0	1.0	1.0	1.0	1.0	1.0	1.0	1.0	1.0
Location	BLK1	GW1	GW2	GW3	GW4	GW5	GW6	SW1	SW2
Field Blank				Field Dup. of CKY19			Field Dup. of CKY16		
CQOL	COMPOUND								
0.05	alpha-BHC								
0.05	beta-BHC								
0.05	delta-BHC								
0.05	*gamma-BHC (Lindane)					0.0039	J		
0.05	*Heptachlor							0.0099	J
0.05	Aldrin								
0.05	Heptachlor Epoxide								
0.05	Endosulfan I								
0.10	Dieldrin								
0.10	4,4'-DDE								
0.10	*Endrin								
0.10	Endosulfan II								
0.10	4,4'-DDD								
0.10	Endosulfan Sulfate								
0.10	4,4'-DDT								
0.50	*Methoxychlor								
0.10	Endrin Ketone								
0.50	*alpha-Chlordane								
0.50	*gamma-Chlordane								
1.0	*Toxaphene								
0.50	*Aroclor-1016								
0.50	*Aroclor-1221								
0.50	*Aroclor-1232								
0.50	*Aroclor-1242								
0.50	*Aroclor-1248								
1.0	*Aroclor-1254								
1.0	*Aroclor-1260								

CQOL = Contract Required Quantitation Limit

ACTION LEVEL EXISTS

BEG NARRATIVE FOR CODE DEFINITION

revised 07/9

DATA SUMMARY FORM: PESTICIDES AND PCPs

Site Name: HOFFMAN LANDFILLASR #: 18347 Sampling Date(s): 6/23/92WATER SAMPLES
(μ g/L)To calculate sample quantitation limit:
(CRQL * Dilution Factor)

Sample No. Dilution Factor Location	CKY24 1.0 SW3	CKY25 1.0 SW4	CKY26 1.0 LTI												
CRTL	COMPOUND														
0.05	alpha-BHC														
0.05	beta-BHC														
0.05	delta-BHC														
0.05	*gamma-BHC (Lindane)														
0.05	*Heptachlor														
0.05	Aldrin														
0.05	Heptachlor Epoxide														
0.05	Endosulfan I														
0.10	Dieldrin														
0.10	4,4'-DDE														
0.10	*Endrin														
0.10	Endosulfan II														
0.10	4,4'-DDD														
0.10	Endosulfan Sulfate					0.0082 J									
0.10	4,4'-DDT														
0.50	*Methoxychlor														
0.10	Endrin Ketone						0.019 B								
0.50	*alpha-Chlordane														
0.50	*gamma-Chlordane														
1.0	*Toxaphene														
0.50	*Aroclor-1016														
0.50	*Aroclor-1221														
0.50	*Aroclor-1232														
0.50	*Aroclor-1242														
0.50	*Aroclor-1248														
1.0	*Aroclor-1254														
1.0	*Aroclor-1260														

CRTL = Contract Required Quantitation Limit

ACTION LEVEL EXISTS

SEE NARRATIVE FOR CODE DEFINITION

DATA SUMMARY FORM: PESTICIDES AND PCB'S

Site Name: HOFFMAN LANDFILL

SOIL SAMPLES
($\mu\text{g}/\text{Kg}$)

Case #: 18347 Sampling Date(s): 6/23/92

To calculate sample quantitation limit:

(CRQL * Dilution Factor) / ((100 - % moisture)/100)

CROL	COMPOUND	CKY 28	CKY 29	CKY 30	CKY 31	CKY 32	CKY 33	CKY 34	CKY 35	CKY 36
		Sample No.	Dilution Factor	X Moisture	Location	1.0	1.0	1.0	1.0	1.0
1.7	alpha-BHC									
1.7	beta-BHC									
1.7	delta-BHC									
1.7	gamma-BHC (Lindane)									
1.7	Heptachlor	0.17 J	0.35 J	0.48 J	0.25 J	0.25 J	0.16 J	0.13 J	0.41 J	0.37 J
1.7	Aldrin									
1.7	Heptachlor Epoxide									
1.7	Endosulfan I									
3.3	Dieldrin			0.072 J		0.20 J	0.25 J			
3.3	4,4'-DDE	0.20 J	0.64 J	0.39 J	0.38 J				0.100 J	
3.3	Endrin		0.47 J				0.075 J	0.59 J	0.27 J	0.51 J
3.3	Endosulfan II							0.31 J	0.22 J	
3.3	4,4'-DDD		0.15 J							
3.3	Endosulfan Sulfate									
3.3	4,4'-DDT									
17	Methoxychlor		0.32 B							
3.3	Endrin Ketone	0.18 J	0.31 J							
1.7	alpha-Chlordane		0.36 J							
1.7	gamma-Chlordane		0.46 J							
170	Toxaphene									
33	Aroclor-1016									
67	Aroclor-1221									
33	Aroclor-1232									
33	Aroclor-1242									
33	Aroclor-1248									
33	Aroclor-1254									
33	Aroclor-1260									
3.3	Endrin Aldehyde									

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DATA SUMMARY FORM: PESTICIDES AND PCB'S

Site Name: HOFFMAN LANDFILL.

Case #: 18347 Sampling Date(s): 6/23/92

SOIL SAMPLES
($\mu\text{g}/\text{Kg}$)To calculate sample quantitation limit:
(CRQL * Dilution Factor) / ((100 - % moisture)/100)

CRQL	COMPOUND	CKY37		CKY38											
		Dilution Factor	% Moisture	Dilution Factor	% Moisture										
1.7	alpha-BHC														
1.7	beta-BHC														
1.7	delta-BHC														
1.7	gamma-BHC (Lindane)	0.12	J												
1.7	Heptachlor			0.094	J										
1.7	Aldrin			0.37	J										
1.7	Heptachlor Epoxide			0.26	J										
1.7	Endosulfan I			0.34	J										
3.3	Dieldrin			0.33	J										
3.3	4,4'-DDF														
3.3	Endrin			0.56	J										
3.3	Endosulfan II														
3.3	4,4'-DDD			1.3	J										
3.3	Endosulfan Sulfate														
3.3	4,4'-DDT														
17	Methoxychlor			0.67	P										
3.3	Endrin Ketone	0.12	J												
1.7	alpha-Chlordane			0.48	J										
1.7	gamma-Chlordane			0.51	J										
170	Toxaphene														
33	Aroclor-1016														
67	Aroclor-1221														
33	Aroclor-1232														
33	Aroclor-1242														
33	Aroclor-1248														
33	Aroclor-1254														
33	Aroclor-1260														
3.3	Endrin Whole Contract Required	Quantitation	Limit												

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OPTIONAL
MATERIAL

Appendix C

**Results as Reported by the Laboratory
for all Target Compounds**

1A
VOLATILE ORGANICS ANALYSIS DATA SHEETOR/2
EPA SAMPLE NO.
*(red)*Lab Name: COMPUCHEM, RTPContract: 68D10083CKY14Lab Code: COMPU Case No.: 18347SAS No.: _____ SDG No.: CKY14Matrix: (soil/water) WATERLab Sample ID: 499425Sample wt/vol: 5.0 (g/mL) MLLab File ID: CN099425A56Level: (low/med) LOWDate Received: 06/24/92

% Moisture: not dec.

Date Analyzed: 06/25/92GC Column: DB624 ID: 0.530 (mm)Dilution Factor: 1.0

Soil Extract Volume: _____ (uL)

Soil Aliquot Volume: _____ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/L</u>	Q
---------	----------	---	---

74-87-3-----	Chloromethane	10	U
74-83-9-----	Bromomethane	10	U
75-01-4-----	Vinyl Chloride	10	U
75-00-3-----	Chloroethane	10	U
75-09-2-----	Methylene Chloride	22	B
67-64-1-----	Acetone	15	B
75-15-0-----	Carbon Disulfide	10	U
75-35-4-----	1,1-Dichloroethene	10	U
75-34-3-----	1,1-Dichloroethane	10	U
540-59-0-----	1,2-Dichloroethene (total)	10	U
67-66-3-----	Chlroform	10	U
107-06-2-----	1,2-Dichloroethane	10	U
78-93-3-----	2-Butanone	10	U
71-55-6-----	1,1,1-Trichloroethane	10	U
56-23-5-----	Carbon Tetrachloride	10	U
75-27-4-----	Bromodichloromethane	10	U
78-87-5-----	1,2-Dichloropropane	10	U
10061-01-5-----	cis-1,3-Dichloropropene	10	U
79-01-6-----	Trichloroethene	10	U
124-48-1-----	Dibromochloromethane	10	U
79-00-5-----	1,1,2-Trichloroethane	10	U
71-43-2-----	Benzene	10	U
10061-02-6-----	Trans-1,3-Dichloropropene	10	U
75-25-2-----	Bromoform	10	U
108-10-1-----	4-Methyl-2-Pentanone	10	U
591-78-6-----	2-Hexanone	10	U
127-18-4-----	Tetrachloroethene	10	U
79-34-5-----	1,1,2,2-Tetrachloroethane	10	U
108-88-3-----	Toluene	10	U
108-90-7-----	Chlorobenzene	10	U
100-41-4-----	Ethylbenzene	10	U
100-42-5-----	Styrene	10	U
1330-20-7-----	Xylene (total)	10	U

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.
(REDACTED)

CKY15

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 18347

SAS No.: _____ SDG No.: CKY14

Matrix: (soil/water) WATER

Lab Sample ID: 499435

Sample wt/vol: 5.0 (g/mL) ML

Lab File ID: CN099435B56

Level: (low/med) LOW

Date Received: 06/24/92

% Moisture: not dec. _____

Date Analyzed: 06/24/92

Column: DB624 ID: 0.530 (mm)

Dilution Factor: 1.0

Soil Extract Volume: _____ (uL)

Soil Aliquot Volume: _____ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) UG/L	Q
---------	----------	--	---

74-87-3-----	Chloromethane	10	U
74-83-9-----	Bromomethane	10	U
75-01-4-----	Vinyl Chloride	10	U
75-00-3-----	Chloroethane	10	U
75-09-2-----	Methylene Chloride	1	BJ
67-64-1-----	Acetone	10	U
75-15-0-----	Carbon Disulfide	10	U
75-35-4-----	1,1-Dichloroethene	10	U
75-34-3-----	1,1-Dichloroethane	10	U
540-59-0-----	1,2-Dichloroethene (total)	10	U
67-66-3-----	Chloroform	10	U
107-06-2-----	1,2-Dichloroethane	10	U
78-93-3-----	2-Butanone	10	U
71-55-6-----	1,1,1-Trichloroethane	10	U
56-23-5-----	Carbon Tetrachloride	10	U
75-27-4-----	Bromodichloromethane	10	U
78-87-5-----	1,2-Dichloropropane	10	U
10061-01-5-----	cis-1,3-Dichloropropene	10	U
79-01-6-----	Trichloroethene	10	U
124-48-1-----	Dibromochloromethane	10	U
79-00-5-----	1,1,2-Trichloroethane	10	U
71-43-2-----	Benzene	10	U
10061-02-6-----	Trans-1,3-Dichloropropene	10	U
75-25-2-----	Bromoform	10	U
108-10-1-----	4-Methyl-2-Pentanone	10	U
591-78-6-----	2-Hexanone	10	U
127-18-4-----	Tetrachloroethene	10	U
79-34-5-----	1,1,2,2-Tetrachloroethane	10	U
108-88-3-----	Toluene	10	U
108-90-7-----	Chlorobenzene	10	U
100-41-4-----	Ethylbenzene	10	U
100-42-5-----	Styrene	10	U
1330-20-7-----	Xylene (total)	10	U

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

CKY16

ORIGINAL
(Rev)

Lab Code: COMPU Case No.: 18347

SAS No.: _____ SDG No.: CKY14

Matrix: (soil/water) WATER

Lab Sample ID: 499436

Sample wt/vol: 5.0 (g/mL) ML

Lab File ID: CN099436B56

Level: (low/med) LOW

Date Received: 06/24/92

% Moisture: not dec. _____

Date Analyzed: 06/24/92

Column: DB624 ID: 0.530 (mm)

Dilution Factor: 1.0

Soil Extract Volume: _____ (uL)

Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

(ug/L or ug/Kg) UG/L

Q

CAS NO.	COMPOUND	UG/L	Q
74-87-3-----	Chloromethane	10	U
74-83-9-----	Bromomethane	10	U
75-01-4-----	Vinyl Chloride	10	U
75-00-3-----	Chloroethane	10	U
75-09-2-----	Methylene Chloride	3	BJ
67-64-1-----	Acetone	15	B
75-15-0-----	Carbon Disulfide	10	U
75-35-4-----	1,1-Dichloroethene	10	U
75-34-3-----	1,1-Dichloroethane	10	U
540-59-0-----	1,2-Dichloroethene (total)	10	U
67-66-3-----	Chloroform	4	J
107-06-2-----	1,2-Dichloroethane	10	U
78-93-3-----	2-Butanone	10	U
71-55-6-----	1,1,1-Trichloroethane	10	U
56-23-5-----	Carbon Tetrachloride	10	U
75-27-4-----	Bromodichloromethane	10	U
78-87-5-----	1,2-Dichloropropane	10	U
10061-01-5-----	cis-1,3-Dichloropropene	10	U
79-01-6-----	Trichloroethene	10	U
124-48-1-----	Dibromochloromethane	10	U
79-00-5-----	1,1,2-Trichloroethane	10	U
71-43-2-----	Benzene	10	U
10061-02-6-----	Trans-1,3-Dichloropropene	10	U
75-25-2-----	Bromoform	10	U
108-10-1-----	4-Methyl-2-Pentanone	10	U
591-78-6-----	2-Hexanone	10	U
127-18-4-----	Tetrachloroethene	10	U
79-34-5-----	1,1,2,2-Tetrachloroethane	10	U
108-88-3-----	Toluene	10	U
108-90-7-----	Chlorobenzene	10	U
100-41-4-----	Ethylbenzene	10	U
100-42-5-----	Styrene	10	U
1330-20-7-----	Xylene (total)	10	U

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CKY17
(Ready)

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 18347

SAS No.: _____ SDG No.: CKY14

Matrix: (soil/water) WATER

Lab Sample ID: 499437

Sample wt/vol: 5.0 (g/mL) ML

Lab File ID: CN099437A56

Level: (low/med) LOW

Date Received: 06/24/92

% Moisture: not dec. _____

Date Analyzed: 06/25/92

Column: DB624 ID: 0.530 (mm)

Dilution Factor: 1.0

Soil Extract Volume: _____ (uL)

Soil Aliquot Volume: _____ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/L</u>	Q
---------	----------	---	---

74-87-3-----	Chloromethane	10	U
74-83-9-----	Bromomethane	10	U
75-01-4-----	Vinyl Chloride	10	U
75-00-3-----	Chloroethane	10	U
75-09-2-----	Methylene Chloride	4	BJ
67-64-1-----	Acetone	9	BJ
75-15-0-----	Carbon Disulfide	10	U
75-35-4-----	1,1-Dichloroethene	10	U
75-34-3-----	1,1-Dichloroethane	10	U
540-59-0-----	1,2-Dichloroethene (total)	10	U
67-66-3-----	Chloroform	10	U
107-06-2-----	1,2-Dichloroethane	10	U
78-93-3-----	2-Butanone	10	U
71-55-6-----	1,1,1-Trichloroethane	10	U
56-23-5-----	Carbon Tetrachloride	10	U
75-27-4-----	Bromodichloromethane	10	U
78-87-5-----	1,2-Dichloropropane	10	U
10061-01-5-----	cis-1,3-Dichloropropene	10	U
79-01-6-----	Trichloroethene	10	U
124-48-1-----	Dibromochloromethane	10	U
79-00-5-----	1,1,2-Trichloroethane	10	U
71-43-2-----	Benzene	10	U
10061-02-6-----	Trans-1,3-Dichloropropene	10	U
75-25-2-----	Bromoform	10	U
108-10-1-----	4-Methyl-2-Pentanone	10	U
591-78-6-----	2-Hexanone	10	U
127-18-4-----	Tetrachloroethene	10	U
79-34-5-----	1,1,2,2-Tetrachloroethane	10	U
108-88-3-----	Toluene	10	U
108-90-7-----	Chlorobenzene	10	U
100-41-4-----	Ethylbenzene	10	U
100-42-5-----	Styrene	10	U
1330-20-7-----	Xylene (total)	10	U

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

ORIGINAL
(Red)

CKY18

Lab Name: COMPUCHEM RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 18347

SAS No.: _____ SDG No.: CKY14

Matrix: (soil/water) WATER

Lab Sample ID: 499439

Sample wt/vol: 5.0 (g/mL) ML

Lab File ID: CR099439C56

Level: (low/med) LOW

Date Received: 06/24/92

% Moisture: not dec. _____

Date Analyzed: 06/26/92

Column: DB624 ID: 0.530 (mm)

Dilution Factor: 1.0

Soil Extract Volume: _____ (uL)

Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

(ug/L or ug/Kg) UG/L

Q

CAS NO.	COMPOUND	UG/L	Q
74-87-3-----	Chloromethane	10	U
74-83-9-----	Bromomethane	10	U
75-01-4-----	Vinyl Chloride	10	U
75-00-3-----	Chlroethane	10	U
75-09-2-----	Methylene Chloride	6	BJ
67-64-1-----	Acetone	12	B
75-15-0-----	Carbon Disulfide	10	U
75-35-4-----	1,1-Dichloroethene	10	U
75-34-3-----	1,1-Dichloroethane	10	U
540-59-0-----	1,2-Dichloroethene (total)	10	U
67-66-3-----	Chloroform	10	U
107-06-2-----	1,2-Dichloroethane	10	U
78-93-3-----	2-Butanone	10	U
71-55-6-----	1,1,1-Trichloroethane	10	U
56-23-5-----	Carbon Tetrachloride	10	U
75-27-4-----	Bromodichloromethane	10	U
78-87-5-----	1,2-Dichloropropane	10	U
10061-01-5-----	cis-1,3-Dichloropropene	10	U
79-01-6-----	Trichloroethene	10	U
124-48-1-----	Dibromochloromethane	10	U
79-00-5-----	1,1,2-Trichloroethane	10	U
71-43-2-----	Benzene	10	U
10061-02-6-----	Trans-1,3-Dichloropropene	10	U
75-25-2-----	Bromoform	10	U
108-10-1-----	4-Methyl-2-Pentanone	10	U
591-78-6-----	2-Hexanone	10	U
127-18-4-----	Tetrachloroethene	10	U
79-34-5-----	1,1,2,2-Tetrachloroethane	10	U
108-88-3-----	Toluene	10	U
108-90-7-----	Chlorobenzene	10	U
100-41-4-----	Ethylbenzene	10	U
100-42-5-----	Styrene	10	U
1330-20-7-----	Xylene (total)	10	U

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

ORIGINAL
(Red)

CKY19

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 18347

SAS No.: _____ SDG No.: CKY14

Matrix: (soil/water) WATER

Lab Sample ID: 499440

Sample wt/vol: 5.0 (g/mL) ML

Lab File ID: CN099440A56

Level: (low/med) LOW

Date Received: 06/24/92

% Moisture: not dec. _____

Date Analyzed: 06/25/92

Column: DB624 ID: 0.530 (mm)

Dilution Factor: 1.0

Soil Extract Volume: _____ (uL)

Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

(ug/L or ug/Kg) UG/L

Q

CAS NO.	COMPOUND	UG/L	Q
74-87-3-----	Chloromethane	10	U
74-83-9-----	Bromomethane	10	U
75-01-4-----	Vinyl Chloride	10	U
75-00-3-----	Chloroethane	10	U
75-09-2-----	Methylene Chloride	6	BJ
67-64-1-----	Acetone	15	B
75-15-0-----	Carbon Disulfide	10	U
75-35-4-----	1,1-Dichloroethene	10	U
75-34-3-----	1,1-Dichloroethane	10	U
540-59-0-----	1,2-Dichloroethene (total)	10	U
67-66-3-----	Chloroform	4	J
107-06-2-----	1,2-Dichloroethane	10	U
78-93-3-----	2-Butanone	10	U
71-55-6-----	1,1,1-Trichloroethane	10	U
56-23-5-----	Carbon Tetrachloride	10	U
75-27-4-----	Bromodichloromethane	10	U
78-87-5-----	1,2-Dichloropropane	10	U
10061-01-5-----	cis-1,3-Dichloropropene	10	U
79-01-6-----	Trichloroethene	10	U
124-48-1-----	Dibromochloromethane	10	U
79-00-5-----	1,1,2-Trichloroethane	10	U
71-43-2-----	Benzene	10	U
10061-02-6-----	Trans-1,3-Dichloropropene	10	U
75-25-2-----	Bromoform	10	U
108-10-1-----	4-Methyl-2-Pentanone	10	U
591-78-6-----	2-Hexanone	10	U
127-18-4-----	Tetrachloroethene	10	U
79-34-5-----	1,1,2,2-Tetrachloroethane	10	U
108-88-3-----	Toluene	10	U
108-90-7-----	Chlorobenzene	10	U
100-41-4-----	Ethylbenzene	10	U
100-42-5-----	Styrene	10	U
1330-20-7-----	Xylene (total)	10	U

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CKY20 ORIGINAL

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 18347

SAS No.: _____ SDG No.: CKY14

Matrix: (soil/water) WATER

Lab Sample ID: 499441

Sample wt/vol: 5.0 (g/mL) ML

Lab File ID: CN099441A56

Level: (low/med) LOW

Date Received: 06/24/92

% Moisture: not dec. _____

Date Analyzed: 06/25/92

Column: DB624 ID: 0.530 (mm)

Dilution Factor: 1.0

Soil Extract Volume: _____ (uL)

Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

(ug/L or ug/Kg) UG/L

Q

CAS NO.	COMPOUND	10	U
74-87-3-----	Chloromethane	10	U
74-83-9-----	Bromomethane	10	U
75-01-4-----	Vinyl Chloride	2	J
75-00-3-----	Chloroethane	10	U
75-09-2-----	Methylene Chloride	3	EJ
67-64-1-----	Acetone	11	B
75-15-0-----	Carbon Disulfide	10	U
75-35-4-----	1,1-Dichloroethene	10	U
75-34-3-----	1,1-Dichloroethane	10	U
540-59-0-----	1,2-Dichloroethene (total)	10	U
67-66-3-----	Chloroform	10	U
107-06-2-----	1,2-Dichloroethane	10	U
78-93-3-----	2-Butanone	10	U
71-55-6-----	1,1,1-Trichloroethane	10	U
56-23-5-----	Carbon Tetrachloride	10	U
75-27-4-----	Bromodichloromethane	10	U
78-87-5-----	1,2-Dichloropropane	10	U
10061-01-5-----	cis-1,3-Dichloropropene	10	U
79-01-6-----	Trichloroethene	10	U
124-48-1-----	Dibromochloromethane	10	U
79-00-5-----	1,1,2-Trichloroethane	10	U
71-43-2-----	Benzene	10	U
10061-02-6-----	Trans-1,3-Dichloropropene	10	U
75-25-2-----	Bromoform	10	U
108-10-1-----	4-Methyl-2-Pentanone	10	U
591-78-6-----	2-Hexanone	10	U
127-18-4-----	Tetrachloroethene	10	U
79-34-5-----	1,1,2,2-Tetrachloroethane	10	U
108-88-3-----	Toluene	10	U
108-90-7-----	Chlorobenzene	10	U
100-41-4-----	Ethylbenzene	10	U
100-42-5-----	Styrene	10	U
1330-20-7-----	Xylene (total)	10	U

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

ORIG.

(Red)

CKY22

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 18347

SAS No.: _____ SDG No.: CKY14

Matrix: (soil/water) WATER

Lab Sample ID: 499442

Sample wt/vol: 5.0 (g/mL) ML

Lab File ID: CN099442A56

Level: (low/med) LOW

Date Received: 06/24/92

% Moisture: not dec. _____

Date Analyzed: 06/25/92

Column: DB624 ID: 0.530 (mm)

Dilution Factor: 1.0

Soil Extract Volume: _____ (uL)

Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

(ug/L or ug/Kg) UG/L

Q

CAS NO.	COMPOUND	UG/L	Q
74-87-3-----	Chloromethane	10	U
74-83-9-----	Bromomethane	10	U
75-01-4-----	Vinyl Chloride	10	U
75-00-3-----	Chloroethane	10	U
75-09-2-----	Methylene Chloride	5	BJ
67-64-1-----	Acetone	12	B
75-15-0-----	Carbon Disulfide	1	J
75-35-4-----	1,1-Dichloroethene	10	U
75-34-3-----	1,1-Dichloroethane	10	U
540-59-0-----	1,2-Dichloroethene (total)	10	U
67-66-3-----	Chloroform	10	U
107-06-2-----	1,2-Dichloroethane	10	U
78-93-3-----	2-Butanone	10	U
71-55-6-----	1,1,1-Trichloroethane	10	U
56-23-5-----	Carbon Tetrachloride	10	U
75-27-4-----	Bromodichloromethane	10	U
78-87-5-----	1,2-Dichloropropane	10	U
10061-01-5-----	cis-1,3-Dichloropropene	10	U
79-01-6-----	Trichloroethene	10	U
124-48-1-----	Dibromochloromethane	10	U
79-00-5-----	1,1,2-Trichloroethane	10	U
71-43-2-----	Benzene	10	U
10061-02-6-----	Trans-1,3-Dichloropropene	10	U
75-25-2-----	Bromoform	10	U
108-10-1-----	4-Methyl-2-Pentanone	10	U
591-78-6-----	2-Hexanone	10	U
127-18-4-----	Tetrachloroethene	10	U
79-34-5-----	1,1,2,2-Tetrachloroethane	10	U
108-88-3-----	Toluene	10	U
108-90-7-----	Chlorobenzene	10	U
100-41-4-----	Ethylbenzene	10	U
100-42-5-----	Styrene	10	U
1330-20-7-----	Xylene (total)	10	U

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

ORIGINAL
(Red)

CKY23

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 18347

SAS No.: _____ SDG No.: CKY14

Matrix: (soil/water) WATER

Lab Sample ID: 499444

Sample wt/vol: 5.0 (g/mL) ML

Lab File ID: CN099444A56

Level: (low/med) LOW

Date Received: 06/24/92

% Moisture: not dec. _____

Date Analyzed: 06/25/92

Column: DB624 ID: 0.530 (mm)

Dilution Factor: 1.0

Soil Extract Volume: _____ (uL)

Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

(ug/L or ug/Kg) UG/L

Q

CAS NO.	COMPOUND	UG/L	Q
74-87-3-----	Chloromethane	10	U
74-83-9-----	Bromomethane	10	U
75-01-4-----	Vinyl Chloride	10	U
75-00-3-----	Chloroethane	10	U
75-09-2-----	Methylene Chloride	5	BJ
67-64-1-----	Acetone	12	B
75-15-0-----	Carbon Disulfide	10	U
75-35-4-----	1,1-Dichloroethene	10	U
75-34-3-----	1,1-Dichloroethane	10	U
540-59-0-----	1,2-Dichloroethene (total)	10	U
67-66-3-----	Chloroform	10	U
107-06-2-----	1,2-Dichloroethane	10	U
78-93-3-----	2-Butanone	10	U
71-55-6-----	1,1,1-Trichloroethane	10	U
56-23-5-----	Carbon Tetrachloride	10	U
75-27-4-----	Bromodichloromethane	10	U
78-87-5-----	1,2-Dichloropropane	10	U
10061-01-5-----	cis-1,3-Dichloropropene	10	U
79-01-6-----	Trichloroethene	10	U
124-48-1-----	Dibromochloromethane	10	U
79-00-5-----	1,1,2-Trichloroethane	10	U
71-43-2-----	Benzene	10	U
10061-02-6-----	Trans-1,3-Dichloropropene	10	U
75-25-2-----	Bromoform	10	U
108-10-1-----	4-Methyl-2-Pentanone	10	U
591-78-6-----	2-Hexanone	10	U
127-18-4-----	Tetrachloroethene	10	U
79-34-5-----	1,1,2,2-Tetrachloroethane	10	U
108-88-3-----	Toluene	10	U
108-90-7-----	Chlorobenzene	10	U
100-41-4-----	Ethylbenzene	10	U
100-42-5-----	Styrene	10	U
1330-20-7-----	Xylene (total)	10	U

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

(REG/NAI)
(Rev.)

CKY24

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 18347

SAS No.: _____ SDG No.: CKY14

Matrix: (soil/water) WATER

Lab Sample ID: 499445

Sample wt/vol: 5.0 (g/mL) ML

Lab File ID: CN099445A56

Level: (low/med) LOW

Date Received: 06/24/92

% Moisture: not dec. _____

Date Analyzed: 06/25/92

Column: DB624 ID: 0.530 (mm)

Dilution Factor: 1.0

Soil Extract Volume: _____ (uL)

Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

(ug/L or ug/Kg) UG/L

Q

CAS NO.	COMPOUND	Q
74-87-3-----	Chloromethane	10 U
74-83-9-----	Bromomethane	10 U
75-01-4-----	Vinyl Chloride	10 U
75-00-3-----	Chlороethane	10 U
75-09-2-----	Methylene Chloride	10 B
67-64-1-----	Acetone	17 B
75-15-0-----	Carbon Disulfide	10 U
75-35-4-----	1,1-Dichloroethene	10 U
75-34-3-----	1,1-Dichloroethane	10 U
540-59-0-----	1,2-Dichloroethene (total)	10 U
67-66-3-----	Chloroform	10 U
107-06-2-----	1,2-Dichloroethane	10 U
78-93-3-----	2-Butanone	10 U
71-55-6-----	1,1,1-Trichloroethane	10 U
56-23-5-----	Carbon Tetrachloride	10 U
75-27-4-----	Bromodichloromethane	10 U
78-87-5-----	1,2-Dichloropropane	10 U
10061-01-5-----	cis-1,3-Dichloropropene	10 U
79-01-6-----	Trichloroethene	10 U
124-48-1-----	Dibromochloromethane	10 U
79-00-5-----	1,1,2-Trichloroethane	10 U
71-43-2-----	Benzene	10 U
10061-02-6-----	Trans-1,3-Dichloropropene	10 U
75-25-2-----	Bromoform	10 U
108-10-1-----	4-Methyl-2-Pentanone	10 U
591-78-6-----	2-Hexanone	10 U
127-18-4-----	Tetrachloroethene	10 U
79-34-5-----	1,1,2,2-Tetrachloroethane	10 U
108-88-3-----	Toluene	10 U
108-90-7-----	Chlorobenzene	10 U
100-41-4-----	Ethylbenzene	10 U
100-42-5-----	Styrene	10 U
1330-20-7-----	Xylene (total)	10 U

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

(REGD)
CKY25

Lab Name: COMPUCHEM RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 18347

SAS No.: _____ SDG No.: CKY14

Matrix: (soil/water) WATER

Lab Sample ID: 499447

Sample wt/vol: 5.0 (g/mL) ML

Lab File ID: CN099447B56

Level: (low/med) LOW

Date Received: 06/24/92

% Moisture: not dec. _____

Date Analyzed: 06/25/92

Column: DB624 ID: 0.530 (mm)

Dilution Factor: 1.0

Soil Extract Volume: _____ (uL)

Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

(ug/L or ug/Kg) UG/L

Q

<u>74-87-3-----Chloromethane</u>	<u>10</u>	<u>U</u>
<u>74-83-9-----Bromomethane</u>	<u>10</u>	<u>U</u>
<u>75-01-4-----Vinyl Chloride</u>	<u>10</u>	<u>U</u>
<u>75-00-3-----Chloroethane</u>	<u>10</u>	<u>U</u>
<u>75-09-2-----Methylene Chloride</u>	<u>3</u>	<u>BJ</u>
<u>67-64-1-----Acetone</u>	<u>11</u>	<u>B</u>
<u>75-15-0-----Carbon Disulfide</u>	<u>10</u>	<u>U</u>
<u>75-35-4-----1,1-Dichloroethene</u>	<u>10</u>	<u>U</u>
<u>75-34-3-----1,1-Dichloroethane</u>	<u>10</u>	<u>U</u>
<u>540-59-0-----1,2-Dichloroethene (total)</u>	<u>10</u>	<u>U</u>
<u>67-66-3-----Chloroform</u>	<u>10</u>	<u>U</u>
<u>107-06-2-----1,2-Dichloroethane</u>	<u>10</u>	<u>U</u>
<u>78-93-3-----2-Butanone</u>	<u>10</u>	<u>U</u>
<u>71-55-6-----1,1,1-Trichloroethane</u>	<u>10</u>	<u>U</u>
<u>56-23-5-----Carbon Tetrachloride</u>	<u>10</u>	<u>U</u>
<u>75-27-4-----Bromodichloromethane</u>	<u>10</u>	<u>U</u>
<u>78-87-5-----1,2-Dichloropropane</u>	<u>10</u>	<u>U</u>
<u>10061-01-5-----cis-1,3-Dichloropropene</u>	<u>10</u>	<u>U</u>
<u>79-01-6-----Trichloroethene</u>	<u>10</u>	<u>U</u>
<u>124-48-1-----Dibromochloromethane</u>	<u>10</u>	<u>U</u>
<u>79-00-5-----1,1,2-Trichloroethane</u>	<u>10</u>	<u>U</u>
<u>71-43-2-----Benzene</u>	<u>10</u>	<u>U</u>
<u>10061-02-6-----Trans-1,3-Dichloropropene</u>	<u>10</u>	<u>U</u>
<u>75-25-2-----Bromoform</u>	<u>10</u>	<u>U</u>
<u>108-10-1-----4-Methyl-2-Pentanone</u>	<u>10</u>	<u>U</u>
<u>591-78-6-----2-Hexanone</u>	<u>10</u>	<u>U</u>
<u>127-18-4-----Tetrachloroethene</u>	<u>10</u>	<u>U</u>
<u>79-34-5-----1,1,2,2-Tetrachloroethane</u>	<u>10</u>	<u>U</u>
<u>108-88-3-----Toluene</u>	<u>10</u>	<u>U</u>
<u>108-90-7-----Chlorobenzene</u>	<u>10</u>	<u>U</u>
<u>100-41-4-----Ethylbenzene</u>	<u>10</u>	<u>U</u>
<u>100-42-5-----Styrene</u>	<u>10</u>	<u>U</u>
<u>1330-20-7-----Xylene (total)</u>	<u>10</u>	<u>U</u>

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

ORIGINAL
(Red)
EPA SAMPLE NO.

CKY26

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 18347 SAS No.: _____ SDG No.: CKY14

Matrix: (soil/water) WATER

Lab Sample ID: 499449

Sample wt/vol: 5.0 (g/mL) ML

Lab File ID: CR099449C56

Level: (low/med) LOW

Date Received: 06/24/92

% Moisture: not dec. _____

Date Analyzed: 06/26/92

GC Column: DB624 ID: 0.530 (mm)

Dilution Factor: 1.0

Soil Extract Volume: _____ (uL)

Soil Aliquot Volume: _____ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/L</u>	Q
---------	----------	---	---

74-87-3-----	Chloromethane	10	U
74-83-9-----	Bromomethane	10	U
75-01-4-----	Vinyl Chloride	10	U
75-00-3-----	Chloroethane	10	U
75-09-2-----	Methylene Chloride	5	BJ
67-64-1-----	Acetone	9	BJ
75-15-0-----	Carbon Disulfide	10	U
75-35-4-----	1,1-Dichloroethene	10	U
75-34-3-----	1,1-Dichloroethane	10	U
540-59-0-----	1,2-Dichloroethene (total)	10	U
67-66-3-----	Chlcroform	10	U
107-06-2-----	1,2-Dichloroethane	10	U
78-93-3-----	2-Butanone	10	U
71-55-6-----	1,1,1-Trichloroethane	10	U
56-23-5-----	Carbon Tetrachloride	10	U
75-27-4-----	Bromodichloromethane	10	U
78-87-5-----	1,2-Dichloropropane	10	U
10061-01-5-----	cis-1,3-Dichloropropene	10	U
79-01-6-----	Trichloroethene	10	U
124-48-1-----	Dibromochloromethane	10	U
79-00-5-----	1,1,2-Trichloroethane	10	U
71-43-2-----	Benzene	10	U
10061-02-6-----	Trans-1,3-Dichloropropene	10	U
75-25-2-----	Bromoform	10	U
108-10-1-----	4-Methyl-2-Pentanone	10	U
591-78-6-----	2-Hexanone	10	U
127-18-4-----	Tetrachloroethene	10	U
79-34-5-----	1,1,2,2-Tetrachloroethane	10	U
108-88-3-----	Toluene	10	U
108-90-7-----	Chlorobenzene	10	U
100-41-4-----	Ethylbenzene	10	U
100-42-5-----	Styrene	10	U
1330-20-7-----	Xylene (total)	10	U

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

CKY28

Lab Code: COMPU Case No.: 18347

SAS No.: _____

SDG No.: **CKY28**

Matrix: (soil/water) SOIL

Lab Sample ID: **499452**

Sample wt/vol: 5.0 (g/mL) G

Lab File ID: **GH099452B54**

Level: (low/med) LOW

Date Received: 06/24/92

% Moisture: not dec. 19

Date Analyzed: 06/24/92

GC Column: DB-624 ID: 0.530 (mm)

Dilution Factor: 1.0

Soil Extract Volume: _____ (uL)

Soil Aliquot Volume: _____ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/KG</u>	Q
74-87-3-----	Chloromethane	12	U
74-83-9-----	Bromomethane	12	U
75-01-4-----	Vinyl Chloride	12	U
75-00-3-----	Chloroethane	12	U
75-09-2-----	Methylene Chloride	30	B
67-64-1-----	Acetone	20	B
75-15-0-----	Carbon Disulfide	12	U
75-35-4-----	1,1-Dichloroethene	12	U
75-34-3-----	1,1-Dichloroethane	12	U
540-59-0-----	1,2-Dichloroethene (total)	12	U
67-66-3-----	Chloroform	12	U
107-06-2-----	1,2-Dichloroethane	12	U
78-93-3-----	2-Butanone	12	U
71-55-6-----	1,1,1-Trichloroethane	12	U
56-23-5-----	Carbon Tetrachloride	12	U
75-27-4-----	Bromodichloromethane	12	U
78-87-5-----	1,2-Dichloroproppane	12	U
10061-01-5-----	cis-1,3-Dichloropropene	12	U
79-01-6-----	Trichloroethene	12	U
124-48-1-----	Dibromochloromethane	12	U
79-00-5-----	1,1,2-Trichloroethane	12	U
71-43-2-----	Benzene	12	U
10061-02-6-----	Trans-1,3-Dichloropropene	12	U
75-25-2-----	Bromoform	12	U
108-10-1-----	4-Methyl-2-Pentanone	12	U
591-78-6-----	2-Hexanone	12	U
127-18-4-----	Tetrachloroethene	12	U
79-34-5-----	1,1,2,2-Tetrachloroethane	12	U
108-88-3-----	Toluene	1	J
108-90-7-----	Chlorobenzene	12	U
100-41-4-----	Ethylbenzene	12	U
100-42-5-----	Styrene	12	U
1330-20-7-----	Xylene (total)	12	U

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 18347

SAS No.: _____ SDG No.: CKY28

Matrix: (soil/water) SOIL

Lab Sample ID: 499459

Sample wt/vol: 5.0 (g/mL) G

Lab File ID: GH099459B54

Level: (low/med) LOW

Date Received: 06/24/92

% Moisture: not dec. 17

Date Analyzed: 06/24/92

GC Column: DB-624 ID: 0.530 (mm)

Dilution Factor: 1.0

Soil Extract Volume: _____ (uL)

Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

(ug/L or ug/Kg) UG/KG

Q

74-87-3-----	Chloromethane	12	U
74-83-9-----	Bromomethane	12	U
75-01-4-----	Vinyl Chloride	12	U
75-00-3-----	Chloroethane	12	U
75-09-2-----	Methylene Chloride	78	B
67-64-1-----	Acetone	32	B
75-15-0-----	Carbon Disulfide	12	U
75-35-4-----	1,1-Dichloroethene	12	U
75-34-3-----	1,1-Dichloroethane	12	U
540-59-0-----	1,2-Dichloroethene (total)	12	U
67-66-3-----	Chloroform	12	U
107-06-2-----	1,2-Dichloroethane	12	U
78-93-3-----	2-Butanone	12	U
71-55-6-----	1,1,1-Trichloroethane	12	U
56-23-5-----	Carbon Tetrachloride	12	U
75-27-4-----	Bromodichloromethane	12	U
78-87-5-----	1,2-Dichloropropane	12	U
10061-01-5-----	cis-1,3-Dichloropropene	12	U
79-01-6-----	Trichloroethene	12	U
124-48-1-----	Dibromochloromethane	12	U
79-00-5-----	1,1,2-Trichloroethane	12	U
71-43-2-----	Benzene	12	U
10061-02-6-----	Trans-1,3-Dichloropropene	12	U
75-25-2-----	Bromoform	12	U
108-10-1-----	4-Methyl-2-Pentanone	12	U
591-78-6-----	2-Hexanone	12	U
127-18-4-----	Tetrachloroethene	12	U
79-34-5-----	1,1,2,2-Tetrachloroethane	12	U
108-88-3-----	Toluene	12	U
108-90-7-----	Chlorobenzene	12	U
100-41-4-----	Ethylbenzene	12	U
100-42-5-----	Styrene	12	U
1330-20-7-----	Xylene (total)	12	U

LA
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.
(Red)

CKY30

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU

Case No.: 18347

SAS No.: _____

SDG No.: CKY28

Matrix: (soil/water) SOIL

Lab Sample ID: 499463

Sample wt/vol: 5.0 (g/mL) G

Lab File ID: GH099463B54

Level: (low/med) LOW

Date Received: 06/24/92

% Moisture: not dec. 40

Date Analyzed: 06/24/92

GC Column: DB-624 ID: 0.530 (mm)

Dilution Factor: 1.0

Soil Extract Volume: _____ (uL)

Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

(ug/L or ug/Kg) UG/KG

Q

CAS NO.	COMPOUND	17	U
74-87-3-----	Chloromethane	17	U
74-83-9-----	Bromomethane	17	U
75-01-4-----	Vinyl Chloride	17	U
75-00-3-----	Chlороethane	17	U
75-09-2-----	Methylene Chloride	84	B
67-64-1-----	Acetone	42	B
75-15-0-----	Carbon Disulfide	17	U
75-35-4-----	1,1-Dichloroethene	17	U
75-34-3-----	1,1-Dichloroethane	17	U
540-59-0-----	1,2-Dichloroethene (total)	17	U
67-66-3-----	Chloroform	17	U
107-06-2-----	1,2-Dichloroethane	17	U
78-93-3-----	2-Butanone	17	U
71-55-6-----	1,1,1-Trichloroethane	17	U
56-23-5-----	Carbon Tetrachloride	17	U
75-27-4-----	Bromodichloromethane	17	U
78-87-5-----	1,2-Dichloropropane	17	U
10061-01-5-----	cis-1,3-Dichloropropene	17	U
79-01-6-----	Trichloroethene	17	U
124-48-1-----	Dibromochloromethane	17	U
79-00-5-----	1,1,2-Trichloroethane	17	U
71-43-2-----	Benzene	17	U
10061-02-6-----	Trans-1,3-Dichloropropene	17	U
75-25-2-----	Bromoform	17	U
108-10-1-----	4-Methyl-2-Pentanone	17	U
591-78-6-----	2-Hexanone	17	U
127-18-4-----	Tetrachloroethene	17	U
79-34-5-----	1,1,2,2-Tetrachloroethane	17	U
108-88-3-----	Toluene	17	U
108-90-7-----	Chlorobenzene	17	U
100-41-4-----	Ethylbenzene	17	U
100-42-5-----	Styrene	17	U
1330-20-7-----	Xylene (total)	17	U

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CKY31

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 18347

SAS No.: _____ SDG No.: CKY28

Matrix: (soil/water) SOIL

Lab Sample ID: 499464

Sample wt/vol: 5.0 (g/mL) G

Lab File ID: GH099464B54

Level: (low/med) LOW

Date Received: 06/24/92

% Moisture: not dec. 23

Date Analyzed: 06/24/92

Column: DB-624 ID: 0.530 (mm)

Dilution Factor: 1.0

Soil Extract Volume: _____ (uL)

Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

(ug/L or ug/Kg) UG/KG

Q

CAS NO.	COMPOUND			
74-87-3-----	Chloromethane	13	U	
74-83-9-----	Bromomethane	13	U	
75-01-4-----	Vinyl Chloride	13	U	
75-00-3-----	Chloroethane	13	U	
75-09-2-----	Methylene Chloride	100	B	
67-64-1-----	Acetone	110	B	
75-15-0-----	Carbon Disulfide	13	U	
75-35-4-----	1,1-Dichloroethene	13	U	
75-34-3-----	1,1-Dichloroethane	13	U	
540-59-0-----	1,2-Dichloroethene (total)	13	U	
67-66-3-----	Chloroform	13	U	
107-06-2-----	1,2-Dichloroethane	13	U	
78-93-3-----	2-Butanone	13	U	
71-55-6-----	1,1,1-Trichloroethane	13	U	
56-23-5-----	Carbon Tetrachloride	13	U	
75-27-4-----	Bromo dichloromethane	13	U	
78-87-5-----	1,2-Dichloropropane	13	U	
10061-01-5-----	cis-1,3-Dichloropropene	13	U	
79-01-6-----	Trichloroethene	13	U	
124-48-1-----	Dibromochloromethane	13	U	
79-00-5-----	1,1,2-Trichloroethane	13	U	
71-43-2-----	Benzene	13	U	
10061-02-6-----	Trans-1,3-Dichloropropene	13	U	
75-25-2-----	Bromoform	13	U	
108-10-1-----	4-Methyl-2-Pentanone	13	U	
591-78-6-----	2-Hexanone	13	U	
127-18-4-----	Tetrachloroethene	13	U	
79-34-5-----	1,1,2,2-Tetrachloroethane	13	U	
108-88-3-----	Toluene	13	U	
108-90-7-----	Chlorobenzene	13	U	
100-41-4-----	Ethylbenzene	13	U	
100-42-5-----	Styrene	13	U	
1330-20-7-----	Xylene (total)	13	U	

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

Og
(Recd)

CKY32

Lab Name: COMPUCHEM RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 18347

SAS No.: _____ SDG No.: CKY28

Matrix: (soil/water) SOIL

Lab Sample ID: 499465

Sample wt/vol: 5.0 (g/mL) G

Lab File ID: GR099465A54

Level: (low/med) LOW

Date Received: 06/24/92

% Moisture: not dec. 18

Date Analyzed: 06/25/92

Column: DB-624 ID: 0.530 (mm)

Dilution Factor: 1.0

Soil Extract Volume: _____ (uL)

Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

(ug/L or ug/Kg) UG/KG

Q

CAS NO.	COMPOUND			
74-87-3-----	Chloromethane	12	U	
74-83-9-----	Bromomethane	12	U	
75-01-4-----	Vinyl Chloride	12	U	
75-00-3-----	Chloroethane	12	U	
75-09-2-----	Methylene Chloride	33	B	
67-64-1-----	Acetone	23	B	
75-15-0-----	Carbon Disulfide	12	U	
75-35-4-----	1,1-Dichloroethene	12	U	
75-34-3-----	1,1-Dichloroethane	12	U	
540-59-0-----	1,2-Dichloroethene (total)	12	U	
67-66-3-----	Chloroform	12	U	
107-06-2-----	1,2-Dichloroethane	12	U	
78-93-3-----	2-Butanone	12	U	
71-55-6-----	1,1,1-Trichloroethane	12	U	
56-23-5-----	Carbon Tetrachloride	12	U	
75-27-4-----	Bromodichloromethane	12	U	
78-87-5-----	1,2-Dichloropropane	12	U	
10061-01-5-----	cis-1,3-Dichloropropene	12	U	
79-01-6-----	Trichloroethene	12	U	
124-48-1-----	Dibromochloromethane	12	U	
79-00-5-----	1,1,2-Trichloroethane	12	U	
71-43-2-----	Benzene	12	U	
10061-02-6-----	Trans-1,3-Dichloropropene	12	U	
75-25-2-----	Bromoform	12	U	
108-10-1-----	4-Methyl-2-Pentanone	12	U	
591-78-6-----	2-Hexanone	12	U	
127-18-4-----	Tetrachloroethene	12	U	
79-34-5-----	1,1,2,2-Tetrachloroethane	12	U	
108-88-3-----	Toluene	12	U	
108-90-7-----	Chlorobenzene	12	U	
100-41-4-----	Ethylbenzene	12	U	
100-42-5-----	Styrene	12	U	
1330-20-7-----	Xylene (total)	12	U	

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

ORIGINAL
(med)

CKY33

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU

Case No.: 18347

SAS No.: _____

SDG No.: CKY28

Matrix: (soil/water) SOIL

Lab Sample ID: 499466

Sample wt/vol: 5.0 (g/mL) G

Lab File ID: GR099466A54

Level: (low/med) LOW

Date Received: 06/24/92

% Moisture: not dec. 11

Date Analyzed: 06/25/92

GC Column: DB-624 ID: 0.530 (mm)

Dilution Factor: 1.0

Soil Extract Volume: _____ (uL)

Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

(ug/L or ug/Kg) UG/KG

Q

74-87-3-----	Chloromethane	11	U
74-83-9-----	Bromomethane	11	U
75-01-4-----	Vinyl Chloride	11	U
75-00-3-----	Chloroethane	11	U
75-09-2-----	Methylene Chloride	31	B
67-64-1-----	Acetone	23	B
75-15-0-----	Carbon Disulfide	11	U
75-35-4-----	1,1-Dichloroethene	11	U
75-34-3-----	1,1-Dichloroethane	11	U
540-59-0-----	1,2-Dichloroethene (total)	11	U
67-66-3-----	Chloroform	11	U
107-06-2-----	1,2-Dichloroethane	11	U
78-93-3-----	2-Butanone	11	U
71-55-6-----	1,1,1-Trichloroethane	11	U
56-23-5-----	Carbon Tetrachloride	11	U
75-27-4-----	Bromodichloromethane	11	U
78-87-5-----	1,2-Dichloropropane	11	U
10061-01-5-----	cis-1,3-Dichloropropene	11	U
79-01-6-----	Trichloroethene	11	U
124-48-1-----	Dibromochloromethane	11	U
79-00-5-----	1,1,2-Trichloroethane	11	U
71-43-2-----	Benzene	11	U
10061-02-6-----	Trans-1,3-Dichloropropene	11	U
75-25-2-----	Bromoform	11	U
108-10-1-----	4-Methyl-2-Pentanone	11	U
591-78-6-----	2-Hexanone	11	U
127-18-4-----	Tetrachloroethene	11	U
79-34-5-----	1,1,2,2-Tetrachloroethane	11	U
108-88-3-----	Toluene	11	U
108-90-7-----	Chlorobenzene	11	U
100-41-4-----	Ethylbenzene	11	U
100-42-5-----	Styrene	11	U
1330-20-7-----	Xylene (total)	11	U

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

ORIGINAL
(Red)

CKY34

Lab Name: COMPUCHEM RTPContract: 68D10083Lab Code: COMPU Case No.: 18347SAS No.: _____ SDG No.: CKY28Matrix: (soil/water) SOILLab Sample ID: 499467Sample wt/vol: 5.0 (g/mL) GLab File ID: GH099467B54Level: (low/med) LOWDate Received: 06/24/92% Moisture: not dec. 13Date Analyzed: 06/24/92GC Column: DB-624 ID: 0.530 (mm)Dilution Factor: 1.0

Soil Extract Volume: _____ (uL)

Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

(ug/L or ug/Kg) UG/KG

Q

CAS NO.	COMPOUND	Q
74-87-3-----	Chloromethane	11 U
74-83-9-----	Bromomethane	11 U
75-01-4-----	Vinyl Chloride	11 U
75-00-3-----	Chloroethane	11 U
75-09-2-----	Methylene Chloride	48 B
67-64-1-----	Acetone	27 B
75-15-0-----	Carbon Disulfide	11 U
75-35-4-----	1,1-Dichloroethene	11 U
75-34-3-----	1,1-Dichloroethane	11 U
540-59-0-----	1,2-Dichloroethene (total)	11 U
67-66-3-----	Chloroform	11 U
107-06-2-----	1,2-Dichloroethane	11 U
78-93-3-----	2-Butanone	11 U
71-55-6-----	1,1,1-Trichloroethane	11 U
56-23-5-----	Carbon Tetrachloride	11 U
75-27-4-----	Bromodichloromethane	11 U
78-87-5-----	1,2-Dichloropropane	11 U
10061-01-5-----	cis-1,3-Dichloropropene	11 U
79-01-6-----	Trichloroethene	11 U
124-48-1-----	Dibromochloromethane	11 U
79-00-5-----	1,1,2-Trichloroethane	11 U
71-43-2-----	Benzene	11 U
10061-02-6-----	Trans-1,3-Dichloropropene	11 U
75-25-2-----	Bromoform	11 U
108-10-1-----	4-Methyl-2-Pentanone	11 U
591-78-6-----	2-Hexanone	11 U
127-18-4-----	Tetrachloroethene	11 U
79-34-5-----	1,1,2,2-Tetrachloroethane	11 U
108-88-3-----	Toluene	11 U
108-90-7-----	Chlorobenzene	11 U
100-41-4-----	Ethylbenzene	11 U
100-42-5-----	Styrene	11 U
1330-20-7-----	Xylene (total)	11 U

1A
VOLATILE ORGANICS ANALYSIS DATA SHEETEPA SAMPLE NO:
(826)

CKY35

Lab Name: COMPUCHEM, RTPContract: 68D10083Lab Code: COMPU Case No.: 18347 SAS No.: _____SDG No.: CKY28Matrix: (soil/water) SOILLab Sample ID: 499468Sample wt/vol: 5.0 (g/mL) GLab File ID: GH099468B54Level: (low/med) LOWDate Received: 06/24/92% Moisture: not dec. 3Date Analyzed: 06/25/92GC Column: DB-624 ID: 0.530 (mm)Dilution Factor: 1.0

Soil Extract Volume: _____ (uL)

Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

(ug/L or ug/Kg) UG/KG

Q

CAS NO.	COMPOUND	Q
74-87-3-----	Chloromethane	10 U
74-83-9-----	Bromomethane	10 U
75-01-4-----	Vinyl Chloride	10 U
75-00-3-----	Chloroethane	10 U
75-09-2-----	Methylene Chloride	87 B
67-64-1-----	Acetone	27 B
75-15-0-----	Carbon Disulfide	10 U
75-35-4-----	1,1-Dichloroethene	10 U
75-34-3-----	1,1-Dichloroethane	10 U
540-59-0-----	1,2-Dichloroethene (total)	10 U
67-66-3-----	Chloroform	10 U
107-06-2-----	1,2-Dichloroethane	10 U
78-93-3-----	2-Butanone	10 U
71-55-6-----	1,1,1-Trichloroethane	10 U
56-23-5-----	Carbon Tetrachloride	10 U
75-27-4-----	Bromodichloromethane	10 U
78-87-5-----	1,2-Dichloropropane	10 U
10061-01-5-----	cis-1,3-Dichloropropene	10 U
79-01-6-----	Trichloroethene	10 U
124-48-1-----	Dibromochloromethane	10 U
79-00-5-----	1,1,2-Trichloroethane	10 U
71-43-2-----	Benzene	10 U
10061-02-6-----	Trans-1,3-Dichloropropene	10 U
75-25-2-----	Bromoform	10 U
108-10-1-----	4-Methyl-2-Pentanone	10 U
591-78-6-----	2-Hexanone	10 U
127-18-4-----	Tetrachloroethene	10 U
79-34-5-----	1,1,2,2-Tetrachloroethane	10 U
108-88-3-----	Toluene	10 U
108-90-7-----	Chlorobenzene	10 U
100-41-4-----	Ethylbenzene	10 U
100-42-5-----	Styrene	10 U
1330-20-7-----	Xylene (total)	10 U

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.
(med)

Lab Name: COMPUCHEM.RTP

Contract: 68D10083

CKY36

Lab Code: COMPU

Case No.: 18347

SAS No.: _____

SDG No.: CKY28

Matrix: (soil/water) SOIL

Lab Sample ID: 499470

Sample wt/vol: 5.0 (g/mL) G

Lab File ID: GR099470A54

Level: (low/med) LOW

Date Received: 06/24/92

Moisture: not dec. 10

Date Analyzed: 06/25/92

GC Column: DB-624 ID: 0.530 (mm)

Dilution Factor: 1.0

Soil Extract Volume: _____ (uL)

CONCENTRATION UNITS:

(ug/L or ug/Kg) UG/KG

Q

CAS NO.	COMPOUND	Q	
74-87-3-----	Chloromethane	11	U
74-83-9-----	Bromomethane	11	U
75-01-4-----	Vinyl Chloride	11	U
75-00-3-----	Chloroethane	11	U
75-09-2-----	Methylene Chloride	43	B
67-64-1-----	Acetone	62	B
75-15-0-----	Carbon Disulfide	11	U
75-35-4-----	1,1-Dichloroethene	11	U
75-34-3-----	1,1-Dichloroethane	11	U
540-59-0-----	1,2-Dichloroethene (total)	11	U
67-66-3-----	Chloroform	11	U
107-06-2-----	1,2-Dichloroethane	11	U
78-93-3-----	2-Butanone	11	U
71-55-6-----	1,1,1-Trichloroethane	11	U
56-23-5-----	Carbon Tetrachloride	11	U
75-27-4-----	Bromodichloromethane	11	U
78-87-5-----	1,2-Dichloropropane	11	U
10061-01-5-----	cis-1,3-Dichloropropene	11	U
79-01-6-----	Trichloroethene	11	U
124-48-1-----	Dibromochloromethane	11	U
79-00-5-----	1,1,2-Trichloroethane	11	U
71-43-2-----	Benzene	11	U
10061-02-6-----	Trans-1,3-Dichloropropene	11	U
75-25-2-----	Bromoform	11	U
108-10-1-----	4-Methyl-2-Pentanone	11	U
591-78-6-----	2-Hexanone	11	U
127-18-4-----	Tetrachloroethene	11	U
79-34-5-----	1,1,2,2-Tetrachloroethane	11	U
108-88-3-----	Toluene	11	U
108-90-7-----	Chlorobenzene	11	U
100-41-4-----	Ethylbenzene	11	U
100-42-5-----	Styrene	11	U
1330-20-7-----	Xylene (total)	11	U

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

ORIGINAL
(Rev)

CKY37

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 18347

SAS No.: _____ SDG No.: CKY28

Matrix: (soil/water) SOIL

Lab Sample ID: 499471

Sample wt/vol: 5.0 (g/mL) G

Lab File ID: GH099471B54

Level: (low/med) LOW

Date Received: 06/24/92

% Moisture: not dec. 8

Date Analyzed: 06/25/92

Column: DB-624 ID: 0.530 (mm)

Dilution Factor: 1.0

Soil Extract Volume: _____ (uL)

Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

CAS NO.	COMPOUND	(ug/L or ug/Kg) UG/KG	Q
---------	----------	-----------------------	---

74-87-3-----	Chloromethane	11	U
74-83-9-----	Bromomethane	11	U
75-01-4-----	Vinyl Chloride	11	U
75-00-3-----	Chloroethane	11	U
75-09-2-----	Methylene Chloride	33	B
67-64-1-----	Acetone	21	B
75-15-0-----	Carbon Disulfide	11	U
75-35-4-----	1,1-Dichloroethene	11	U
75-34-3-----	1,1-Dichloroethane	11	U
540-59-0-----	1,2-Dichloroethene (total)	11	U
67-66-3-----	Chloroform	11	U
107-06-2-----	1,2-Dichloroethane	11	U
78-93-3-----	2-Butanone	11	U
71-55-6-----	1,1,1-Trichloroethane	11	U
56-23-5-----	Carbon Tetrachloride	11	U
75-27-4-----	Bromodichloromethane	11	U
78-87-5-----	1,2-Dichloropropane	11	U
10061-01-5-----	cis-1,3-Dichloropropene	11	U
79-01-6-----	Trichloroethene	11	U
124-48-1-----	Dibromochloromethane	11	U
79-00-5-----	1,1,2-Trichloroethane	11	U
71-43-2-----	Benzene	11	U
10061-02-6-----	Trans-1,3-Dichloropropene	11	U
75-25-2-----	Bromoform	11	U
108-10-1-----	4-Methyl-2-Pentanone	11	U
591-78-6-----	2-Hexanone	11	U
127-18-4-----	Tetrachloroethene	11	U
79-34-5-----	1,1,2,2-Tetrachloroethane	11	U
108-88-3-----	Toluene	11	U
108-90-7-----	Chlorobenzene	11	U
100-41-4-----	Ethylbenzene	11	U
100-42-5-----	Styrene	11	U
1330-20-7-----	Xylene (total)	11	U

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

ORIGINAL
(Red)

CKY38

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 18347

SAS No.: _____ SDG No.: CKY28

Matrix: (soil/water) SOIL

Lab Sample ID: 499473

Sample wt/vol: 5.0 (g/mL) G

Lab File ID: GH099473C54

Level: (low/med) LOW

Date Received: 06/24/92

% Moisture: not dec. 12

Date Analyzed: 06/25/92

GC Column: DB-624 ID: 0.530 (mm)

Dilution Factor: 1.0

Soil Extract Volume: _____ (uL)

Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

(ug/L or ug/Kg) UG/KG

Q

74-87-3-----	Chloromethane	11	U
74-83-9-----	Bromomethane	11	U
75-01-4-----	Vinyl Chloride	11	U
75-00-3-----	Chloroethane	11	U
75-09-2-----	Methylene Chloride	98	B
67-64-1-----	Acetone	35	B
75-15-0-----	Carbon Disulfide	11	U
75-35-4-----	1,1-Dichloroethene	11	U
75-34-3-----	1,1-Dichloroethane	11	U
540-59-0-----	1,2-Dichloroethene (total)	11	U
67-66-3-----	Chloroform	11	U
107-06-2-----	1,2-Dichloroethane	11	U
78-93-3-----	2-Butanone	11	U
71-55-6-----	1,1,1-Trichloroethane	11	U
56-23-5-----	Carbon Tetrachloride	11	U
75-27-4-----	Bromodichloromethane	11	U
78-87-5-----	1,2-Dichloropropane	11	U
10061-01-5-----	cis-1,3-Dichloropropene	11	U
79-01-6-----	Trichloroethene	11	U
124-48-1-----	Dibromochloromethane	11	U
79-00-5-----	1,1,2-Trichloroethane	11	U
71-43-2-----	Benzene	11	U
10061-02-6-----	Trans-1,3-Dichloropropene	11	U
75-25-2-----	Bromoform	11	U
108-10-1-----	4-Methyl-2-Pentanone	11	U
591-78-6-----	2-Hexanone	11	U
127-18-4-----	Tetrachloroethene	11	U
79-34-5-----	1,1,2,2-Tetrachloroethane	11	U
108-88-3-----	Toluene	11	U
108-90-7-----	Chlorobenzene	11	U
100-41-4-----	Ethylbenzene	11	U
100-42-5-----	Styrene	11	U
1330-20-7-----	Xylene (total)	11	U

1B
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

ORIGINAL
CKY14 (Red)Lab Name: COMPUCHEM, RTPContract: 68D10083Lab Code: COMPU Case No.: 18347SAS No.: _____ SDG No.: CKY14Matrix: (soil/water) WATERLab Sample ID: 499425Sample wt/vol: 1000 (g/mL) MLLab File ID: GH099425B02Level: (low/med) LOWDate Received: 06/24/92

% Moisture: _____ decanted: (Y/N) _____

Date Extracted: 06/26/92Concentrated Extract Volume: 1000 (uL)Date Analyzed: 06/29/92Injection Volume: 2.0(uL)Dilution Factor: 1.0GPC Cleanup: (Y/N) N pH: _____

CONCENTRATION UNITS:

(ug/L or ug/Kg) UG/LQ

CAS NO.	COMPOUND	10	U
108-95-2	Phenol	10	U
111-44-4	bis(2-Chloroethyl)Ether	10	U
95-57-8	2-Chlorophenol	10	U
541-73-1	1,3-Dichlorobenzene	10	U
106-46-7	1,4-Dichlorobenzene	10	U
95-50-1	1,2-Dichlorobenzene	10	U
95-48-7	2-Methylphenol	10	U
108-60-1	2,2'-Oxybis(1-Chloropropane)	10	U
106-44-5	4-Methylphenol	10	U
621-64-7	N-Nitroso-Di-n-Propylamine	10	U
67-72-1	Hexachloroethane	10	U
98-95-3	Nitrobenzene	10	U
78-59-1	Isophorone	10	U
88-75-5	2-Nitrophenol	10	U
105-67-9	2,4-Dimethylphenol	10	U
111-91-1	bis(2-Chloroethoxy)Methane	10	U
120-83-2	2,4-Dichlorophenol	10	U
120-82-1	1,2,4-Trichlorobenzene	10	U
91-20-3	Naphthalene	10	U
106-47-8	4-Chloroaniline	10	U
87-68-3	Hexachlorobutadiene	10	U
59-50-7	4-Chloro-3-Methylphenol	10	U
91-57-6	2-Methylnaphthalene	10	U
77-47-4	Hexachlorocyclopentadiene	10	U
88-06-2	2,4,6-Trichlorophenol	10	U
95-95-4	2,4,5-Trichlorophenol	25	U
91-58-7	2-Chloronaphthalene	10	U
88-74-4	2-Nitroaniline	25	U
131-11-3	Dimethyl Phthalate	10	U
208-96-8	Acenaphthylene	10	U
606-20-2	2,6-Dinitrotoluene	10	U
99-09-2	3-Nitroaniline	25	U
83-32-9	Acenaphthene	10	U

SEMOVOLATILE ORGANICS ANALYSIS DATA SHEET

CKY14

Lab Name: COMPUCHEM, RTP Contract: 68D10083

Lab Code: COMPU Case No.: 18347 SAS No.: _____ SDG No.: CKY14

Matrix: (soil/water) WATER Lab Sample ID: 499425

Sample wt/vol: 1000 (g/mL) ML Lab File ID: GH099425B02

Level: (low/med) LOW Date Received: 06/24/92

% Moisture: _____ decanted: (Y/N) _____ Date Extracted: 06/26/92

Concentrated Extract Volume: 1000 (uL) Date Analyzed: 06/29/92

Injection Volume: 2.0 (uL) Dilution Factor: 1.0

GPC Cleanup: (Y/N) N pH: _____

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/L</u>	Q
51-28-5-----	2,4-Dinitrophenol	25	U
100-02-7-----	4-Nitrophenol	25	U
132-64-9-----	Dibenzofuran	10	U
121-14-2-----	2,4-Dinitrotoluene	10	U
84-66-2-----	Diethylphthalate	10	U
7005-72-3-----	4-Chlorophenyl-phenylether	10	U
86-73-7-----	Fluorene	10	U
100-01-6-----	4-Nitroaniline	25	U
534-52-1-----	4,6-Dinitro-2-Methylphenol	25	U
86-30-6-----	N-Nitrosodiphenylamine (1)	10	U
101-55-3-----	4-Bromophenyl-phenylether	10	U
118-74-1-----	Hexachlorobenzene	10	U
87-86-5-----	Pentachlorophenol	25	U
85-01-8-----	Phenanthrene	10	U
120-12-7-----	Anthracene	10	U
86-74-8-----	Carbazole	10	U
84-74-2-----	Di-n-Butylphthalate	10	U
206-44-0-----	Fluoranthene	10	U
129-00-0-----	Pyrene	10	U
85-68-7-----	Butylbenzylphthalate	10	U
91-94-1-----	3,3'-Dichlorobenzidine	10	U
56-55-3-----	Benzo(a)Anthracene	10	U
218-01-9-----	Chrysene	10	U
117-81-7-----	bis(2-Ethylhexyl)Phthalate	2	BJ
117-84-0-----	Di-n-Octyl Phthalate	10	U
205-99-2-----	Benzo(b)Fluoranthene	10	U
207-08-9-----	Benzo(k)Fluoranthene	10	U
50-32-8-----	Benzo(a)Pyrene	10	U
193-39-5-----	Indeno(1,2,3-cd)Pyrene	10	U
53-70-3-----	Dibenz(a,h)Anthracene	10	U
191-24-2-----	Benzo(g,h,i)Perylene	10	U

(1) - Cannot be separated from Diphenylamine

1B
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.
(Red)
GHAI

CKY15

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 18347

SAS No.: _____ SDG No.: CKY14

Matrix: (soil/water) WATER

Lab Sample ID: 499435

Sample wt/vol: 1000 (g/mL) ML

Lab File ID: GH099435B02

Level: (low/med) LOW

Date Received: 06/24/92

% Moisture: _____ decanted: (Y/N)

Date Extracted: 06/26/92

Concentrated Extract Volume: 1000 (uL)

Date Analyzed: 06/29/92

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) N pH:

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/L Q

CAS NO.	COMPOUND	10	U
108-95-2-----	Phenol	10	U
111-44-4-----	bis(2-Chloroethyl)Ether	10	U
95-57-8-----	2-Chlorophenol	10	U
541-73-1-----	1,3-Dichlorobenzene	10	U
106-46-7-----	1,4-Dichlorobenzene	10	U
95-50-1-----	1,2-Dichlorobenzene	10	U
95-48-7-----	2-Methylphenol	10	U
108-60-1-----	2,2'-Oxybis(1-Chloropropane)	10	U
106-44-5-----	4-Methylphenol	10	U
621-64-7-----	N-Nitroso-Di-n-Propylamine	10	U
67-72-1-----	Hexachloroethane	10	U
98-95-3-----	Nitrobenzene	10	U
78-59-1-----	Isophorone	10	U
88-75-5-----	2-Nitrophenol	10	U
105-67-9-----	2,4-Dimethylphenol	10	U
111-91-1-----	bis(2-Chloroethoxy)Methane	10	U
120-83-2-----	2,4-Dichlorophenol	10	U
120-82-1-----	1,2,4-Trichlorobenzene	10	U
91-20-3-----	Naphthalene	10	U
106-47-8-----	4-Chloroaniline	10	U
87-68-3-----	Hexachlorobutadiene	10	U
59-50-7-----	4-Chloro-3-Methylphenol	10	U
91-57-6-----	2-Methylnaphthalene	10	U
77-47-4-----	Hexachlorocyclopentadiene	10	U
88-06-2-----	2,4,6-Trichlorophenol	10	U
95-95-4-----	2,4,5-Trichlorophenol	25	U
91-58-7-----	2-Chloronaphthalene	10	U
88-74-4-----	2-Nitroaniline	25	U
131-11-3-----	Dimethyl Phthalate	10	U
208-96-8-----	Acenaphthylene	10	U
606-20-2-----	2,6-Dinitrotoluene	10	U
99-09-2-----	3-Nitroaniline	25	U
83-32-9-----	Acenaphthene	10	U

1C
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEETORIGINAL
EPA SAMPLE NO.Lab Name: COMPUCHEM, RTPContract: 68D10083CKY15Lab Code: COMPUCase No.: 18347

SAS No.: _____

SDG No.: CKY14Matrix: (soil/water) WATERLab Sample ID: 499435Sample wt/vol: 1000 (g/mL) MLLab File ID: GH099435B02Level: (low/med) LOWDate Received: 06/24/92

% Moisture: _____ decanted: (Y/N) _____

Date Extracted: 06/26/92Concentrated Extract Volume: 1000 (uL)Date Analyzed: 06/29/92Injection Volume: 2.0 (uL)Dilution Factor: 1.0GPC Cleanup: (Y/N) N pH: _____CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/L

Q

CAS NO.	COMPOUND	UNITS	Q
51-28-5-----	2,4-Dinitrophenol	25	U
100-02-7-----	4-Nitrophenol	25	U
132-64-9-----	Dibenzofuran	10	U
121-14-2-----	2,4-Dinitrotoluene	10	U
84-66-2-----	Diethylphthalate	10	U
7005-72-3-----	4-Chlorophenyl-phenylether	10	U
86-73-7-----	Fluorene	10	U
100-01-6-----	4-Nitroaniline	25	U
534-52-1-----	4,6-Dinitro-2-Methylphenol	25	U
86-30-6-----	N-Nitrosodiphenylamine (1)	10	U
101-55-3-----	4-Bromophenyl-phenylether	10	U
118-74-1-----	Hexachlorobenzene	10	U
87-86-5-----	Pentachlorophenol	25	U
85-01-8-----	Phenanthrene	10	U
120-12-7-----	Anthracene	10	U
86-74-8-----	Carbazole	10	U
84-74-2-----	Di-n-Butylphthalate	10	U
206-44-0-----	Fluoranthene	10	U
129-00-0-----	Pyrene	10	U
85-68-7-----	Butylbenzylphthalate	10	U
91-94-1-----	3,3'-Dichlorobenzidine	10	U
56-55-3-----	Benzo(a)Anthracene	10	U
218-01-9-----	Chrysene	10	U
117-81-7-----	bis(2-Ethylhexyl)Phthalate	2	BJ
117-84-0-----	Di-n-Octyl Phthalate	10	U
205-99-2-----	Benzo(b)Fluoranthene	10	U
207-08-9-----	Benzo(k)Fluoranthene	10	U
50-32-8-----	Benzo(a)Pyrene	10	U
193-39-5-----	Indeno(1,2,3-cd)Pyrene	10	U
53-70-3-----	Dibenz(a,h)Anthracene	10	U
191-24-2-----	Benzo(g,h,i)Perylene	10	U

(1) - Cannot be separated from Diphenylamine

1B
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

ORIGINAL
(Rev)

CKY16

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU

Case No.: 18347

SAS No.: _____

SDG No.: CKY14

Matrix: (soil/water) WATER

Lab Sample ID: 499436

Sample wt/vol: 1000 (g/mL) ML

Lab File ID: GH099436B02

Level: (low/med) LOW

Date Received: 06/24/92

% Moisture: _____ decanted: (Y/N) _____

Date Extracted: 06/26/92

Concentrated Extract Volume: 1000 (uL)

Date Analyzed: 06/29/92

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) N pH: _____

CONCENTRATION UNITS:

(ug/L or ug/Kg) UG/L

Q

CAS NO.	COMPOUND	10	U
108-95-2	Phenol	10	U
111-44-4	bis(2-Chloroethyl) Ether	10	U
95-57-8	2-Chlorophenol	10	U
541-73-1	1,3-Dichlorobenzene	10	U
106-46-7	1,4-Dichlorobenzene	10	U
95-50-1	1,2-Dichlorobenzene	10	U
95-48-7	2-Methylphenol	10	U
108-60-1	2,2'-Oxybis(1-Chloropropane)	10	U
106-44-5	4-Methylphenol	10	U
621-64-7	N-Nitroso-Di-n-Propylamine	10	U
67-72-1	Hexachloroethane	10	U
98-95-3	Nitrobenzene	10	U
78-59-1	Isophorone	10	U
88-75-5	2-Nitrophenol	10	U
105-67-9	2,4-Dimethylphenol	10	U
111-91-1	bis(2-Chloroethoxy) Methane	10	U
120-83-2	2,4-Dichlorophenol	10	U
120-82-1	1,2,4-Trichlorobenzene	10	U
91-20-3	Naphthalene	10	U
106-47-8	4-Chloroaniline	10	U
87-68-3	Hexachlorobutadiene	10	U
59-50-7	4-Chloro-3-Methylphenol	10	U
91-57-6	2-Methylnaphthalene	10	U
77-47-4	Hexachlorocyclopentadiene	10	U
88-06-2	2,4,6-Trichlorophenol	10	U
95-95-4	2,4,5-Trichlorophenol	25	U
91-58-7	2-Chloronaphthalene	10	U
88-74-4	2-Nitroaniline	25	U
131-11-3	Dimethyl Phthalate	10	U
208-96-8	Acenaphthylene	10	U
606-20-2	2,6-Dinitrotoluene	10	U
99-09-2	3-Nitroaniline	25	U
83-32-9	Acenaphthene	10	U

1C
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

OR
ORIGINAL
(P&D)
CKY16Lab Name: COMPUCHEM, RTPContract: 68D10083Lab Code: COMPUCase No.: 18347

SAS No.: _____

SDG No.: CKY14Matrix: (soil/water) WATERLab Sample ID: 499436Sample wt/vol: 1000 (g/mL) MLLab File ID: GH099436B02Level: (low/med) LOWDate Received: 06/24/92

% Moisture: _____ decanted: (Y/N) _____

Date Extracted: 06/26/92Concentrated Extract Volume: 1000 (uL)Date Analyzed: 06/29/92Injection Volume: 2.0 (uL)Dilution Factor: 1.0GPC Cleanup: (Y/N) N pH: _____

CONCENTRATION UNITS:

(ug/L or ug/Kg) UG/L

Q

CAS NO.	COMPOUND	25	U
51-28-5-----	2,4-Dinitrophenol	25	U
100-02-7-----	4-Nitrophenol	25	U
132-64-9-----	Dibenzofuran	10	U
121-14-2-----	2,4-Dinitrotoluene	10	U
84-66-2-----	Diethylphthalate	10	U
7005-72-3-----	4-Chlorophenyl-phenylether	10	U
86-73-7-----	Fluorene	10	U
100-01-6-----	4-Nitroaniline	25	U
534-52-1-----	4,6-Dinitro-2-Methylphenol	25	U
86-30-6-----	N-Nitrosodiphenylamine (1)	10	U
101-55-3-----	4-Bromophenyl-phenylether	10	U
118-74-1-----	Hexachlorobenzene	10	U
87-86-5-----	Pentachlorophenol	25	U
85-01-8-----	Phenanthrene	10	U
120-12-7-----	Anthracene	10	U
86-74-8-----	Carbazole	10	U
84-74-2-----	Di-n-Butylphthalate	10	U
206-44-0-----	Fluoranthene	10	U
129-00-0-----	Pyrene	10	U
85-68-7-----	Butylbenzylphthalate	10	U
91-94-1-----	3,3'-Dichlorobenzidine	10	U
56-55-3-----	Benzo(a)Anthracene	10	U
218-01-9-----	Chrysene	10	U
117-81-7-----	bis(2-Ethylhexyl)Phthalate	1	BJ
117-84-0-----	Di-n-Octyl Phthalate	10	U
205-99-2-----	Benzo(b)Fluoranthene	10	U
207-08-9-----	Benzo(k)Fluoranthene	10	U
50-32-8-----	Benzo(a)Pyrene	10	U
193-39-5-----	Indeno(1,2,3-cd)Pyrene	10	U
53-70-3-----	Dibenz(a,h)Anthracene	10	U
191-24-2-----	Benzo(g,h,i)Perylene	10	U

(1) - Cannot be separated from Diphenylamine

1B
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

ORIGINAL
(red)

CKY17

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 18347

SAS No.: _____ SDG No.: CKY14

Matrix: (soil/water) WATER

Lab Sample ID: 499437

Sample wt/vol: 1000 (g/mL) ML

Lab File ID: GH099437B02

Level: (low/med) LOW

Date Received: 06/24/92

% Moisture: _____ decanted: (Y/N) _____

Date Extracted: 06/26/92

Concentrated Extract Volume: 1000 (uL)

Date Analyzed: 06/30/92

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) N pH: _____

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/L Q

CAS NO.	COMPOUND	10	U
108-95-2-----	Phenol	10	U
111-44-4-----	bis(2-Chloroethyl)Ether	10	U
95-57-8-----	2-Chlorophenol	10	U
541-73-1-----	1,3-Dichlorobenzene	10	U
106-46-7-----	1,4-Dichlorobenzene	10	U
95-50-1-----	1,2-Dichlorobenzene	10	U
95-48-7-----	2-Methylphenol	10	U
108-60-1-----	2,2'-Oxybis(1-Chloropropane)	10	U
106-44-5-----	4-Methylphenol	10	U
621-64-7-----	N-Nitroso-Di-n-Propylamine	10	U
67-72-1-----	Hexachloroethane	10	U
98-95-3-----	Nitrobenzene	10	U
78-59-1-----	Isophorone	10	U
88-75-5-----	2-Nitrophenol	10	U
105-67-9-----	2,4-Dimethylphenol	10	U
111-91-1-----	bis(2-Chloroethoxy)Methane	10	U
120-33-2-----	2,4-Dichlorophenol	10	U
120-32-1-----	1,2,4-Trichlorobenzene	10	U
91-20-3-----	Naphthalene	10	U
106-47-8-----	4-Chloroaniline	10	U
87-68-3-----	Hexachlorobutadiene	10	U
59-50-7-----	4-Chloro-3-Methylphenol	10	U
91-57-6-----	2-Methylnaphthalene	10	U
77-47-4-----	Hexachlorocyclopentadiene	10	U
88-06-2-----	2,4,6-Trichlorophenol	10	U
95-95-4-----	2,4,5-Trichlorophenol	25	U
91-58-7-----	2-Chloronaphthalene	10	U
88-74-4-----	2-Nitroaniline	25	U
131-11-3-----	Dimethyl Phthalate	10	U
208-96-8-----	Acenaphthylene	10	U
606-20-2-----	2,6-Dinitrotoluene	10	U
99-09-2-----	3-Nitroaniline	25	U
83-32-9-----	Acenaphthene	10	U

SEMOVOLATILE ORGANICS ANALYSIS DATA SHEET

CKY17

Lab Name: COMPUCHEM, RTP Contract: 68D10083Lab Code: COMPU Case No.: 18347 SAS No.: _____ SDG No.: CKY14Matrix: (soil/water) WATERLab Sample ID: 499437Sample wt/vol: 1000 (g/mL) MLLab File ID: GH099437B02Level: (low/med) LOWDate Received: 06/24/92

% Moisture: _____ decanted: (Y/N) _____

Date Extracted: 06/26/92Concentrated Extract Volume: 1000 (uL)Date Analyzed: 06/30/92Injection Volume: 2.0(uL)Dilution Factor: 1.0GPC Cleanup: (Y/N) N pH: _____CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/L Q

CAS NO.	COMPOUND	25	U
51-28-5-----	2,4-Dinitrophenol	25	U
100-02-7-----	4-Nitrophenol	25	U
132-64-9-----	Dibenzofuran	10	U
121-14-2-----	2,4-Dinitrotoluene	10	U
84-66-2-----	Diethylphthalate	10	U
7005-72-3-----	4-Chlorophenyl-phenylether	10	U
86-73-7-----	Fluorene	10	U
100-01-6-----	4-Nitroaniline	25	U
534-52-1-----	4,6-Dinitro-2-Methylphenol	25	U
86-30-6-----	N-Nitrosodiphenylamine (1)	10	U
101-55-3-----	4-Bromophenyl-phenylether	10	U
118-74-1-----	Hexachlorobenzene	10	U
87-86-5-----	Pentachlorophenol	25	U
85-01-8-----	Phenanthrene	10	U
120-12-7-----	Anthracene	10	U
86-74-8-----	Carbazole	10	U
84-74-2-----	Di-n-Butylphthalate	10	U
206-44-0-----	Fluoranthene	10	U
129-00-0-----	Pyrene	10	U
85-68-7-----	Butylbenzylphthalate	10	U
91-94-1-----	3,3'-Dichlorobenzidine	10	U
56-55-3-----	Benzo(a)Anthracene	10	U
218-01-9-----	Chrysene	10	U
117-81-7-----	bis(2-Ethylhexyl)Phthalate	3	BJ
117-84-0-----	Di-n-Octyl Phthalate	10	U
205-99-2-----	Benzo(b)Fluoranthene	10	U
207-08-9-----	Benzo(k)Fluoranthene	10	U
50-32-8-----	Benzo(a)Pyrene	10	U
193-39-5-----	Indeno(1,2,3-cd) Pyrene	10	U
53-70-3-----	Dibenz(a,h)Anthracene	10	U
191-24-2-----	Benzo(g,h,i)Perylene	10	U

(1) - Cannot be separated from Diphenylamine

1B
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CKY18

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU

Case No.: 18347

SAS No.: _____

SDG No.: CKY14

Matrix: (soil/water) WATER

Lab Sample ID: 499439

Sample wt/vol: 1000 (g/mL) ML

Lab File ID: GH099439B02

Level: (low/med) LOW

Date Received: 06/24/92

% Moisture: _____ decanted: (Y/N) _____

Date Extracted: 06/26/92

Concentrated Extract Volume: 1000 (uL)

Date Analyzed: 06/29/92

Injection Volume: 2.0(uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) N pH: _____

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/L

Q

CAS NO.	COMPOUND	10	U
108-95-2-----	Phenol	10	U
111-44-4-----	bis(2-Chloroethyl)Ether	10	U
95-57-8-----	2-Chlorophenol	10	U
541-73-1-----	1,3-Dichlorobenzene	10	U
106-46-7-----	1,4-Dichlorobenzene	10	U
95-50-1-----	1,2-Dichlorobenzene	10	U
95-48-7-----	2-Methylphenol	10	U
108-60-1-----	2,2'-Oxybis(1-Chloropropane)	10	U
106-44-5-----	4-Methylphenol	10	U
621-64-7-----	N-Nitroso-Di-n-Propylamine	10	U
67-72-1-----	Hexachloroethane	10	U
98-95-3-----	Nitrobenzene	10	U
78-59-1-----	Isophorone	10	U
88-75-5-----	2-Nitrophenol	10	U
105-67-9-----	2,4-Dimethylphenol	10	U
111-91-1-----	bis(2-Chloroethoxy)Methane	10	U
120-83-2-----	2,4-Dichlorophenol	10	U
120-82-1-----	1,2,4-Trichlorobenzene	10	U
91-20-3-----	Naphthalene	10	U
106-47-8-----	4-Chloroaniline	10	U
87-68-3-----	Hexachlorobutadiene	10	U
59-50-7-----	4-Chloro-3-Methylphenol	10	U
91-57-6-----	2-Methylnaphthalene	10	U
77-47-4-----	Hexachlorocyclopentadiene	10	U
88-06-2-----	2,4,6-Trichlorophenol	10	U
95-95-4-----	2,4,5-Trichlorophenol	25	U
91-58-7-----	2-Chloronaphthalene	10	U
88-74-4-----	2-Nitroaniline	25	U
131-11-3-----	Dimethyl Phthalate	10	U
208-96-8-----	Acenaphthylene	10	U
606-20-2-----	2,6-Dinitrotoluene	10	U
99-09-2-----	3-Nitroaniline	25	U
83-32-9-----	Acenaphthene	10	U

1C
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

(P/N)
(D/J)

CKY18

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 18347

SAS No.: _____

SDG No.: CKY14

Matrix: (soil/water) WATER

Lab Sample ID: 499439

Sample wt/vol: 1000 (g/mL) ML

Lab File ID: GH099439B02

Level: (low/med) LOW

Date Received: 06/24/92

% Moisture: _____ decanted: (Y/N)

Date Extracted: 06/26/92

Concentrated Extract Volume: 1000 (uL)

Date Analyzed: 06/29/92

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) N pH: _____

CONCENTRATION UNITS:

(ug/L or ug/Kg) UG/L

Q

CAS NO.	COMPOUND	CONCENTRATION UNITS:	Q
51-28-5-----	2,4-Dinitrophenol	25	U
100-02-7-----	4-Nitrophenol	25	U
132-64-9-----	Dibenzofuran	10	U
121-14-2-----	2,4-Dinitrotoluene	10	U
84-66-2-----	Diethylphthalate	10	U
7005-72-3-----	4-Chlorophenyl-phenylether	10	U
86-73-7-----	Fluorene	10	U
100-01-6-----	4-Nitroaniline	25	U
534-52-1-----	4,6-Dinitro-2-Methylphenol	25	U
86-30-6-----	N-Nitrosodiphenylamine (1)	10	U
101-55-3-----	4-Bromophenyl-phenylether	10	U
118-74-1-----	Hexachlorobenzene	10	U
87-86-5-----	Pentachlorophenol	25	U
85-01-8-----	Phenanthrene	10	U
120-12-7-----	Anthracene	10	U
86-74-8-----	Carbazole	10	U
84-74-2-----	Di-n-Butylphthalate	10	U
206-44-0-----	Fluoranthene	10	U
129-00-0-----	Pyrene	10	U
85-68-7-----	Butylbenzylphthalate	10	U
91-94-1-----	3,3'-Dichlorobenzidine	10	U
56-55-3-----	Benzo(a)Anthracene	10	U
218-01-9-----	Chrysene	10	U
117-81-7-----	bis(2-Ethylhexyl)Phthalate	3	BJ
117-84-0-----	Di-n-Octyl Phthalate	10	U
205-99-2-----	Benzo(b)Fluoranthene	10	U
207-08-9-----	Benzo(k)Fluoranthene	10	U
50-32-8-----	Benzo(a)Pyrene	10	U
193-39-5-----	Indeno(1,2,3-cd) Pyrene	10	U
53-70-3-----	Dibenz(a,h)Anthracene	10	U
191-24-2-----	Benzo(g,h,i)Perylene	10	U

(1) - Cannot be separated from Diphenylamine

1B
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.
(Op/13 Med)

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

CKY19

Lab Code: COMPU

Case No.: 18347

SAS No.: _____

SDG No.: CKY14

Matrix: (soil/water) WATER

Sample wt/vol: 1000 (g/mL) ML

Level: (low/med) LOW

% Moisture: _____ decanted: (Y/N) _____

Concentrated Extract Volume: 1000 (uL)

Injection Volume: 2.0 (uL)

GPC Cleanup: (Y/N) N pH: _____

Lab Sample ID: 499440

Lab File ID: GH099440B02

Date Received: 06/24/92

Date Extracted: 06/26/92

Date Analyzed: 06/30/92

Dilution Factor: 1.0

CAS NO. COMPOUND

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/L

Q

108-95-2-----	Phenol	10	U
111-44-4-----	bis(2-Chloroethyl)Ether	10	U
95-57-8-----	2-Chlorophenol	10	U
541-73-1-----	1,3-Dichlorobenzene	10	U
106-46-7-----	1,4-Dichlorobenzene	10	U
95-50-1-----	1,2-Dichlorobenzene	10	U
95-48-7-----	2-Methylphenol	10	U
108-60-1-----	2,2'-Oxybis(1-Chloropropane)	10	U
106-44-5-----	4-Methylphenol	10	U
621-64-7-----	N-Nitroso-Di-n-Propylamine	10	U
67-72-1-----	Hexachloroethane	10	U
98-95-3-----	Nitrobenzene	10	U
78-59-1-----	Isophorone	10	U
88-75-5-----	2-Nitrophenol	10	U
105-67-9-----	2,4-Dimethylphenol	10	U
111-91-1-----	bis(2-Chloroethoxy)Methane	10	U
120-83-2-----	2,4-Dichlorophenol	10	U
120-82-1-----	1,2,4-Trichlorobenzene	10	U
91-20-3-----	Naphthalene	10	U
106-47-8-----	4-Chloroaniline	10	U
87-68-3-----	Hexachlorobutadiene	10	U
59-50-7-----	4-Chloro-3-Methylphenol	10	U
91-57-6-----	2-Methylnaphthalene	10	U
77-47-4-----	Hexachlorocyclopentadiene	10	U
88-06-2-----	2,4,6-Trichlorophenol	10	U
95-95-4-----	2,4,5-Trichlorophenol	25	U
91-58-7-----	2-Chloronaphthalene	10	U
88-74-4-----	2-Nitroaniline	25	U
131-11-3-----	Dimethyl Phthalate	10	U
208-96-8-----	Acenaphthylene	10	U
606-20-2-----	2,6-Dinitrotoluene	10	U
99-09-2-----	3-Nitroaniline	25	U
83-32-9-----	Acenaphthene	10	U

1C
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEETORIGIN
EPA SAMPLE NO.
(med)

CKY19

Lab Name: COMPUCHEM, RTP Contract: 68D10083Lab Code: COMPU Case No.: 18347 SAS No.: _____ SDG No.: CKY14Matrix: (soil/water) WATERLab Sample ID: 499440Sample wt/vol: 1000 (g/mL) MLLab File ID: GH099440B02Level: (low/med) LOWDate Received: 06/24/92

% Moisture: _____ decanted: (Y/N) _____

Date Extracted: 06/26/92Concentrated Extract Volume: 1000 (uL)Date Analyzed: 06/30/92Injection Volume: 2.0 (uL)Dilution Factor: 1.0GPC Cleanup: (Y/N) N pH: _____CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/L Q

CAS NO.	COMPOUND			
51-28-5-----	2,4-Dinitrophenol	25	U	
100-02-7-----	4-Nitrophenol	25	U	
132-64-9-----	Dibenzofuran	10	U	
121-14-2-----	2,4-Dinitrotoluene	10	U	
84-66-2-----	Diethylphthalate	10	U	
7005-72-3-----	4-Chlorophenyl-phenylether	10	U	
86-73-7-----	Fluorene	10	U	
100-01-6-----	4-Nitroaniline	25	U	
534-52-1-----	4,6-Dinitro-2-Methylphenol	25	U	
86-30-6-----	N-Nitrosodiphenylamine (1)	10	U	
101-55-3-----	4-Bromophenyl-phenylether	10	U	
118-74-1-----	Hexachlorobenzene	10	U	
87-86-5-----	Pentachlorophenol	25	U	
85-01-8-----	Phenanthrene	10	U	
120-12-7-----	Anthracene	10	U	
86-74-8-----	Carbazole	10	U	
84-74-2-----	Di-n-Butylphthalate	10	U	
206-44-0-----	Fluoranthene	10	U	
129-00-0-----	Pyrene	10	U	
85-68-7-----	Butylbenzylphthalate	10	U	
91-94-1-----	3,3'-Dichlorobenzidine	10	U	
56-55-3-----	Benzo(a)Anthracene	10	U	
218-01-9-----	Chrysene	10	U	
117-81-7-----	bis(2-Ethylhexyl)Phthalate	2	BJ	
117-84-0-----	Di-n-Octyl Phthalate	10	U	
205-99-2-----	Benzo(b)Fluoranthene	10	U	
207-08-9-----	Benzo(k)Fluoranthene	10	U	
50-32-8-----	Benzo(a)Pyrene	10	U	
193-39-5-----	Indeno(1,2,3-cd)Pyrene	10	U	
53-70-3-----	Dibenz(a,h)Anthracene	10	U	
191-24-2-----	Benzo(g,h,i)Perylene	10	U	

(1) - Cannot be separated from Diphenylamine

1B
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

ORIGINAL
(Red)

CKY20

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU

Case No.: 18347

SAS No.: _____

SDG No.: CKY14

Matrix: (soil/water) WATER

Lab Sample ID: 499441

Sample wt/vol: 1000 (g/mL) ML

Lab File ID: GH099441B02

Level: (low/med) LOW

Date Received: 06/24/92

% Moisture: _____ decanted: (Y/N)

Date Extracted: 06/26/92

Concentrated Extract Volume: 1000 (uL)

Date Analyzed: 06/30/92

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) N pH:

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/L Q

CAS NO.	COMPOUND	10	U
108-95-2-----	Phenol	10	U
111-44-4-----	bis(2-Chloroethyl)Ether	10	U
95-57-8-----	2-Chlorophenol	10	U
541-73-1-----	1,3-Dichlorobenzene	10	U
106-46-7-----	1,4-Dichlorobenzene	10	U
95-50-1-----	1,2-Dichlorobenzene	10	U
95-48-7-----	2-Methylphenol	10	U
108-60-1-----	2,2'-Oxybis(1-Chloropropane)	10	U
106-44-5-----	4-Methylphenol	10	U
621-64-7-----	N-Nitroso-Di-n-Propylamine	10	U
67-72-1-----	Hexachloroethane	10	U
98-95-3-----	Nitrobenzene	10	U
78-59-1-----	Iscophorone	10	U
88-75-5-----	2-Nitrophenol	10	U
105-67-9-----	2,4-Dimethylphenol	10	U
111-91-1-----	bis(2-Chloroethoxy)Methane	10	U
120-83-2-----	2,4-Dichlorophenol	10	U
120-82-1-----	1,2,4-Trichlorobenzene	10	U
91-20-3-----	Naphthalene	10	U
106-47-8-----	4-Chloroaniline	10	U
87-68-3-----	Hexachlorobutadiene	10	U
59-50-7-----	4-Chloro-3-Methylphenol	10	U
91-57-6-----	2-Methylnaphthalene	10	U
77-47-4-----	Hexachlorocyclopentadiene	10	U
88-06-2-----	2,4,6-Trichlorophenol	10	U
95-95-4-----	2,4,5-Trichlorophenol	25	U
91-58-7-----	2-Chloronaphthalene	10	U
88-74-4-----	2-Nitroaniline	25	U
131-11-3-----	Dimethyl Phthalate	10	U
208-96-8-----	Acenaphthylene	10	U
606-20-2-----	2,6-Dinitrotoluene	10	U
99-09-2-----	3-Nitroaniline	25	U
83-32-9-----	Acenaphthene	10	U

1C
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CKY20

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 18347

SAS No.: _____ SDG No.: CKY14

Matrix: (soil/water) WATER

Lab Sample ID: 499441

Sample wt/vol: 1000 (g/mL) ML

Lab File ID: GH099441B02

Level: (low/med) LOW

Date Received: 06/24/92

% Moisture: _____ decanted: (Y/N) _____

Date Extracted: 06/26/92

Concentrated Extract Volume: 1000 (uL)

Date Analyzed: 06/30/92

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) N pH: _____

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/L

CAS NO.	COMPOUND	Q
51-28-5-----	2,4-Dinitrophenol	25 U
100-02-7-----	4-Nitrophenol	25 U
132-64-9-----	Dibenzofuran	10 U
121-14-2-----	2,4-Dinitrotoluene	10 U
84-66-2-----	Diethylphthalate	10 U
7005-72-3-----	4-Chlorophenyl-phenylether	10 U
86-73-7-----	Fluorene	10 U
100-01-6-----	4-Nitroaniline	25 U
534-52-1-----	4,6-Dinitro-2-Methylphenol	25 U
86-30-6-----	N-Nitrosodiphenylamine (1)	10 U
101-55-3-----	4-Bromophenyl-phenylether	10 U
118-74-1-----	Hexachlorobenzene	10 U
87-86-5-----	Pentachlorophenol	25 U
85-01-8-----	Phenanthrene	10 U
120-12-7-----	Anthracene	10 U
86-74-8-----	Carbazole	10 U
84-74-2-----	Di-n-Butylphthalate	10 U
206-44-0-----	Fluoranthene	10 U
129-00-0-----	Pyrene	10 U
85-68-7-----	Butylbenzylphthalate	10 U
91-94-1-----	3,3'-Dichlorobenzidine	10 U
56-55-3-----	Benzo(a)Anthracene	10 U
218-01-9-----	Chrysene	10 U
117-81-7-----	bis(2-Ethylhexyl)Phthalate	6 BJ
117-84-0-----	Di-n-Octyl Phthalate	10 U
205-99-2-----	Benzo(b)Fluoranthene	10 U
207-08-9-----	Benzo(k)Fluoranthene	10 U
50-32-8-----	Benzo(a)Pyrene	10 U
193-39-5-----	Indeno(1,2,3-cd)Pyrene	10 U
53-70-3-----	Dibenz(a,h)Anthracene	10 U
191-24-2-----	Benzo(g,h,i)Perylene	10 U

(1) - Cannot be separated from Diphenylamine

1B
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.
(REG)

CKY22

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 18347

SAS No.: _____

SDG No.: CKY14

Matrix: (soil/water) WATER

Lab Sample ID: 499442

Sample wt/vol: 1000 (g/mL) ML

Lab File ID: GH099442B02

Level: (low/med) LOW

Date Received: 06/24/92

% Moisture: _____ decanted: (Y/N) _____

Date Extracted: 06/26/92

Concentrated Extract Volume: 1000 (uL)

Date Analyzed: 06/30/92

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) N pH: _____

CONCENTRATION UNITS:

(ug/L or ug/Kg) UG/L

Q

CAS NO.	COMPOUND	2	BJ
108-95-2	Phenol	2	BJ
111-44-4	bis(2-Chloroethyl)Ether	10	U
95-57-8	2-Chlorophenol	10	U
541-73-1	1,3-Dichlorobenzene	10	U
106-46-7	1,4-Dichlorobenzene	10	U
95-50-1	1,2-Dichlorobenzene	10	U
95-48-7	2-Methylphenol	10	U
108-60-1	2,2'-Oxybis(1-Chloropropane)	10	U
106-44-5	4-Methylphenol	10	U
621-64-7	N-Nitroso-Di-n-Propylamine	10	U
67-72-1	Hexachloroethane	10	U
98-95-3	Nitrobenzene	10	U
78-59-1	Isophorone	10	U
88-75-5	2-Nitrophenol	10	U
105-67-9	2,4-Dimethylphenol	10	U
111-91-1	bis(2-Chloroethoxy)Methane	10	U
120-83-2	2,4-Dichlorophenol	10	U
120-82-1	1,2,4-Trichlorobenzene	10	U
91-20-3	Naphthalene	10	U
106-47-8	4-Chloroaniline	10	U
87-68-3	Hexachlorobutadiene	10	U
59-50-7	4-Chloro-3-Methylphenol	10	U
91-57-6	2-Methylnaphthalene	10	U
77-47-4	Hexachlorocyclopentadiene	10	U
88-06-2	2,4,6-Trichlorophenol	10	U
95-95-4	2,4,5-Trichlorophenol	25	U
91-58-7	2-Chloronaphthalene	10	U
88-74-4	2-Nitroaniline	25	U
131-11-3	Dimethyl Phthalate	10	U
208-96-8	Acenaphthylene	10	U
606-20-2	2,6-Dinitrotoluene	10	U
99-09-2	3-Nitroaniline	25	U
83-32-9	Acenaphthene	10	U

SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

(Ref ID)

CKY22

Lab Name: COMPUCHEM, RTPContract: 68D10083Lab Code: COMPUCase No.: 18347

SAS No.: _____

SDG No.: CKY14Matrix: (soil/water) WATERLab Sample ID: 499442Sample wt/vol: 1000 (g/mL) MLLab File ID: GH099442B02Level: (low/med) LOWDate Received: 06/24/92Moisture: _____ decanted: (Y/N) Date Extracted: 06/26/92Concentrated Extract Volume: 1000 (uL)Date Analyzed: 06/30/92Injection Volume: 2.0(uL)Dilution Factor: 1.0GPC Cleanup: (Y/N) N pH: _____CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/L

Q

CAS NO.	COMPOUND	25	U
51-28-5-----	2,4-Dinitrophenol	25	U
100-02-7-----	4-Nitrophenol	25	U
132-64-9-----	Dibenzofuran	10	U
121-14-2-----	2,4-Dinitrotoluene	10	U
84-66-2-----	Diethylphthalate	10	U
7005-72-3-----	4-Chlorophenyl-phenylether	10	U
86-73-7-----	Fluorene	10	U
100-01-6-----	4-Nitroaniline	25	U
534-52-1-----	4,6-Dinitro-2-Methylphenol	25	U
86-30-6-----	N-Nitrosodiphenylamine (1)	10	U
101-55-3-----	4-Bromophenyl-phenylether	10	U
118-74-1-----	Hexachlorobenzene	10	U
87-86-5-----	Pentachlorophenol	25	U
85-01-8-----	Phenanthrene	10	U
120-12-7-----	Anthracene	10	U
86-74-8-----	Carbazole	10	U
84-74-2-----	Di-n-Butylphthalate	10	U
206-44-0-----	Fluoranthene	10	U
129-00-0-----	Pyrene	10	U
85-68-7-----	Butylbenzylphthalate	10	U
91-94-1-----	3,3'-Dichlorobenzidine	10	U
56-55-3-----	Benzo(a)Anthracene	10	U
218-01-9-----	Chrysene	10	U
117-81-7-----	bis(2-Ethylhexyl)Phthalate	6	BJ
117-84-0-----	Di-n-Octyl Phthalate	10	U
205-99-2-----	Benzo(b)Fluoranthene	10	U
207-08-9-----	Benzo(k)Fluoranthene	10	U
50-32-8-----	Benzo(a)Pyrene	10	U
193-39-5-----	Indeno(1,2,3-cd)Pyrene	10	U
53-70-3-----	Dibenz(a,h)Anthracene	10	U
191-24-2-----	Benzo(g,h,i)Perylene	10	U

(1) - Cannot be separated from Diphenylamine

1B
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CKY23

ORIGINAL
(REGD)

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU

Case No.: 18347

SAS No.: _____

SDG No.: CKY14

Matrix: (soil/water) WATER

Lab Sample ID: 499444

Sample wt/vol: 1000 (g/mL) ML

Lab File ID: GH099444B02

Level: (low/med) LOW

Date Received: 06/24/92

% Moisture: _____ decanted: (Y/N)

Date Extracted: 06/26/92

Concentrated Extract Volume: 1000 (uL)

Date Analyzed: 06/30/92

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) N pH:

CONCENTRATION UNITS:

(ug/L or ug/Kg) UG/L

Q

CAS NO.	COMPOUND	10	U
108-95-2-----	Phenol	10	U
111-44-4-----	bis(2-Chloroethyl)Ether	10	U
95-57-8-----	2-Chlorophenol	10	U
541-73-1-----	1,3-Dichlorobenzene	10	U
106-46-7-----	1,4-Dichlorobenzene	10	U
95-50-1-----	1,2-Dichlorobenzene	10	U
95-48-7-----	2-Methylphenol	10	U
108-60-1-----	2,2'-Oxybis(1-Chloropropane)	10	U
106-44-5-----	4-Methylphenol	10	U
621-64-7-----	N-Nitroso-Di-n-Propylamine	10	U
67-72-1-----	Hexachloroethane	10	U
98-95-3-----	Nitrobenzene	10	U
78-59-1-----	Isophorone	10	U
88-75-5-----	2-Nitrophenol	10	U
105-67-9-----	2,4-Dimethylphenol	10	U
111-91-1-----	bis(2-Chloroethoxy)Methane	10	U
120-83-2-----	2,4-Dichlorophenol	10	U
120-82-1-----	1,2,4-Trichlorobenzene	10	U
91-20-3-----	Naphthalene	10	U
106-47-8-----	4-Chloroaniline	10	U
87-68-3-----	Hexachlorobutadiene	10	U
59-50-7-----	4-Chloro-3-Methylphenol	10	U
91-57-6-----	2-Methylnaphthalene	10	U
77-47-4-----	Hexachlorocyclopentadiene	10	U
88-06-2-----	2,4,6-Trichlorophenol	10	U
95-95-4-----	2,4,5-Trichlorophenol	25	U
91-58-7-----	2-Chloronaphthalene	10	U
88-74-4-----	2-Nitroaniline	25	U
131-11-3-----	Dimethyl Phthalate	10	U
208-96-8-----	Acenaphthylene	10	U
606-20-2-----	2,6-Dinitrotoluene	10	U
99-09-2-----	3-Nitroaniline	25	U
83-32-9-----	Acenaphthene	10	U

1C
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

Lab Name: COMPUCHEM.RTPContract: 68D10083CKY20 ORIGINAL
(Rev)Lab Code: COMPUCase No.: 18347

SAS No.: _____

SDG No.: CKY14Matrix: (soil/water) WATERLab Sample ID: 499444Sample wt/vol: 1000 (g/mL) MLLab File ID: GH099444B02Level: (low/med) LOWDate Received: 06/24/92% Moisture: _____ decanted: (Y/N) Date Extracted: 06/26/92Concentrated Extract Volume: 1000 (uL)Date Analyzed: 06/30/92Injection Volume: 2.0 (uL)Dilution Factor: 1.0GPC Cleanup: (Y/N) N pH: _____

CONCENTRATION UNITS:

CAS NO.	COMPOUND	(ug/L or ug/Kg)	UG/L	Q
51-28-5-----	2,4-Dinitrophenol	25	U	
100-02-7-----	4-Nitrophenol	25	U	
132-64-9-----	Dibenzofuran	10	U	
121-14-2-----	2,4-Dinitrotoluene	10	U	
84-66-2-----	Diethylphthalate	10	U	
7005-72-3-----	4-Chlorophenyl-phenylether	10	U	
86-73-7-----	Fluorene	10	U	
100-01-6-----	4-Nitroaniline	25	U	
534-52-1-----	4,6-Dinitro-2-Methylphenol	25	U	
86-30-6-----	N-Nitrosodiphenylamine (1)	10	U	
101-55-3-----	4-Bromophenyl-phenylether	10	U	
118-74-1-----	Hexachlorobenzene	10	U	
87-86-5-----	Pentachlorophenol	25	U	
85-01-8-----	Phenanthrene	10	U	
120-12-7-----	Anthracene	10	U	
86-74-8-----	Carbazole	10	U	
84-74-2-----	Di-n-Butylphthalate	10	U	
206-44-0-----	Fluoranthene	10	U	
129-00-0-----	Pyrene	10	U	
85-68-7-----	Butylbenzylphthalate	10	U	
91-94-1-----	3,3'-Dichlorobenzidine	10	U	
56-55-3-----	Benzo(a)Anthracene	10	U	
218-01-9-----	Chrysene	10	U	
117-81-7-----	bis(2-Ethylhexyl)Phthalate	3	BJ	
117-84-0-----	Di-n-Octyl Phthalate	10	U	
205-99-2-----	Benzo(b)Fluoranthene	10	U	
207-08-9-----	Benzo(k)Fluoranthene	10	U	
50-32-8-----	Benzo(a)Pyrene	10	U	
193-39-5-----	Indeno(1,2,3-cd)Pyrene	10	U	
53-70-3-----	Dibenz(a,h)Anthracene	10	U	
191-24-2-----	Benzo(g,h,i)Perylene	10	U	

(1) - Cannot be separated from Diphenylamine

1B
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

ORIGINAL
(Red)

CKY24

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 18347

SAS No.: _____

SDG No.: CKY14

Matrix: (soil/water) WATER

Lab Sample ID: 499445

Sample wt/vol: 1000 (g/mL) ML

Lab File ID: GH099445A07

Level: (low/med) LOW

Date Received: 06/24/92

% Moisture: _____ decanted: (Y/N) _____

Date Extracted: 06/25/92

Concentrated Extract Volume: 1000 (uL)

Date Analyzed: 06/30/92

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) N pH: _____

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/L Q

CAS NO.	COMPOUND	10	U
108-95-2-----	Phenol	10	U
111-44-4-----	bis(2-Chloroethyl)Ether	10	U
95-57-8-----	2-Chlorophenol	10	U
541-73-1-----	1,3-Dichlorobenzene	10	U
106-46-7-----	1,4-Dichlorobenzene	10	U
95-50-1-----	1,2-Dichlorobenzene	10	U
95-48-7-----	2-Methylphenol	10	U
108-60-1-----	2,2'-Oxybis(1-Chloropropane)	10	U
106-44-5-----	4-Methylphenol	10	U
621-64-7-----	N-Nitroso-Di-n-Propylamine	10	U
67-72-1-----	Hexachloroethane	10	U
98-95-3-----	Nitrobenzene	10	U
78-59-1-----	Iscophorone	10	U
88-75-5-----	2-Nitrophenol	10	U
105-67-9-----	2,4-Dimethylphenol	10	U
111-91-1-----	bis(2-Chloroethoxy)Methane	10	U
120-83-2-----	2,4-Dichlorophenol	10	U
120-82-1-----	1,2,4-Trichlorobenzene	10	U
91-20-3-----	Naphthalene	10	U
106-47-8-----	4-Chloroaniline	10	U
87-68-3-----	Hexachlorobutadiene	10	U
59-50-7-----	4-Chloro-3-Methylphenol	10	U
91-57-6-----	2-Methylnaphthalene	10	U
77-47-4-----	Hexachlorocyclopentadiene	10	U
88-06-2-----	2,4,6-Trichlorophenol	10	U
95-95-4-----	2,4,5-Trichlorophenol	25	U
91-58-7-----	2-Chloronaphthalene	10	U
88-74-4-----	2-Nitroaniline	25	U
131-11-3-----	Dimethyl Phthalate	10	U
208-96-8-----	Acenaphthylene	10	U
606-20-2-----	2,6-Dinitrotoluene	10	U
99-09-2-----	3-Nitroaniline	25	U
83-32-9-----	Acenaphthene	10	U

1C
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

ORIGINAL
CKY24 Red

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU

Case No.: 18347

SAS No.: _____

SDG No.: CKY14

Matrix: (soil/water) WATER

Lab Sample ID: 499445

Sample wt/vol: 1000 (g/mL) ML

Lab File ID: GH099445A07

Level: (low/med) LOW

Date Received: 06/24/92

% Moisture: _____ decanted: (Y/N) _____

Date Extracted: 06/25/92

Concentrated Extract Volume: 1000 (uL)

Date Analyzed: 06/30/92

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) N pH: _____

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/L Q

CAS NO.	COMPOUND	CONCENTRATION UNITS:	Q
51-28-5-----	2,4-Dinitrophenol	25	U
100-02-7-----	4-Nitrophenol	25	U
132-64-9-----	Dibenzofuran	10	U
121-14-2-----	2,4-Dinitrotoluene	10	U
84-66-2-----	Diethylphthalate	10	U
7005-72-3-----	4-Chlorophenyl-phenylether	10	U
86-73-7-----	Fluorene	10	U
100-01-6-----	4-Nitroaniline	25	U
534-52-1-----	4,6-Dinitro-2-Methylphenol	25	U
86-30-6-----	N-Nitrosodiphenylamine (1)	10	U
101-55-3-----	4-Bromophenyl-phenylether	10	U
118-74-1-----	Hexachlorobenzene	10	U
87-86-5-----	Pentachlorophenol	25	U
85-01-8-----	Phenanthrene	10	U
120-12-7-----	Anthracene	10	U
86-74-8-----	Carbazole	10	U
84-74-2-----	Di-n-Butylphthalate	10	U
206-44-0-----	Fluoranthene	10	U
129-00-0-----	Pyrene	10	U
85-68-7-----	Butylbenzylphthalate	10	U
91-94-1-----	3,3'-Dichlorobenzidine	10	U
56-55-3-----	Benzo(a)Anthracene	10	U
218-01-9-----	Chrysene	10	U
117-81-7-----	bis(2-Ethylhexyl)Phthalate	3	BJ
117-84-0-----	Di-n-Octyl Phthalate	10	U
205-99-2-----	Benzo(b)Fluoranthene	10	U
207-08-9-----	Benzo(k)Fluoranthene	10	U
50-32-8-----	Benzo(a)Pyrene	10	U
193-39-5-----	Indeno(1,2,3-cd)Pyrene	10	U
53-70-3-----	Dibenz(a,h)Anthracene	10	U
191-24-2-----	Benzo(g,h,i)Perylene	10	U

(1) - Cannot be separated from Diphenylamine

1B
SEMOVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

OP/CAL

(Reed)

CKY25

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU

Case No.: 18347

SAS No.: _____

SDG No.: CKY14

Matrix: (soil/water) WATER

Lab Sample ID: 499447

Sample wt/vol: 1000 (g/mL) ML

Lab File ID: GH099447A07

Level: (low/med) LOW

Date Received: 06/24/92

% Moisture: _____ decanted: (Y/N) _____

Date Extracted: 06/25/92

Concentrated Extract Volume: 1000 (uL)

Date Analyzed: 06/30/92

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) N pH: _____

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/L

Q

CAS NO.	COMPOUND			
108-95-2	Phenol	10	U	
111-44-4	bis(2-Chloroethyl)Ether	10	U	
95-57-8	2-Chlorophenol	10	U	
541-73-1	1,3-Dichlorobenzene	10	U	
106-46-7	1,4-Dichlorobenzene	10	U	
95-50-1	1,2-Dichlorobenzene	10	U	
95-48-7	2-Methylphenol	10	U	
108-60-1	2,2'-Oxybis(1-Chloropropane)	10	U	
106-44-5	4-Methylphenol	10	U	
621-64-7	N-Nitroso-Di-n-Propylamine	10	U	
67-72-1	Hexachloroethane	10	U	
98-95-3	Nitrobenzene	10	U	
78-59-1	Isophorone	10	U	
88-75-5	2-Nitrophenol	10	U	
105-67-9	2,4-Dimethylphenol	10	U	
111-91-1	bis(2-Chloroethoxy)Methane	10	U	
120-83-2	2,4-Dichlorophenol	10	U	
120-32-1	1,2,4-Trichlorobenzene	10	U	
91-20-3	Naphthalene	10	U	
106-47-8	4-Chloroaniline	10	U	
87-68-3	Hexachlorobutadiene	10	U	
59-50-7	4-Chloro-3-Methylphenol	10	U	
91-57-6	2-Methylnaphthalene	10	U	
77-47-4	Hexachlorocyclopentadiene	10	U	
88-06-2	2,4,6-Trichlorophenol	10	U	
95-95-4	2,4,5-Trichlorophenol	25	U	
91-58-7	2-Chloronaphthalene	10	U	
88-74-4	2-Nitroaniline	25	U	
131-11-3	Dimethyl Phthalate	10	U	
208-96-8	Acenaphthylene	10	U	
606-20-2	2,6-Dinitrotoluene	10	U	
99-09-2	3-Nitroaniline	25	U	
83-32-9	Acenaphthene	10	U	

1C
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

ORIGINAL
EPA SAMPLE NO.

CKY25

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 18347

SAS No.: _____

SDG No.: CKY14

Matrix: (soil/water) WATER

Lab Sample ID: 499447

Sample wt/vol: 1000 (g/mL) ML

Lab File ID: GH099447A07

Level: (low/med) LOW

Date Received: 06/24/92

Moisture: _____ decanted: (Y/N) _____

Date Extracted: 06/25/92

Concentrated Extract Volume: 1000 (uL)

Date Analyzed: 06/30/92

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) N pH: _____

CONCENTRATION UNITS:

(ug/L or ug/Kg) UG/L

Q

CAS NO.	COMPOUND	UG/L	Q
51-28-5-----	2,4-Dinitrophenol	25	U
100-02-7-----	4-Nitrophenol	25	U
132-64-9-----	Dibenzofuran	10	U
121-14-2-----	2,4-Dinitrotoluene	10	U
84-66-2-----	Diethylphthalate	10	U
7005-72-3-----	4-Chlorophenyl-phenylether	10	U
86-73-7-----	Fluorene	10	U
100-01-6-----	4-Nitroaniline	25	U
534-52-1-----	4,6-Dinitro-2-Methylphenol	25	U
86-30-6-----	N-Nitrosodiphenylamine (1)	10	U
101-55-3-----	4-Bromophenyl-phenylether	10	U
118-74-1-----	Hexachlorobenzene	10	U
87-86-5-----	Pentachlorophenol	25	U
85-01-8-----	Phenanthrene	10	U
120-12-7-----	Anthracene	10	U
86-74-8-----	Carbazole	10	U
84-74-2-----	Di-n-Butylphthalate	10	U
206-44-0-----	Fluoranthene	10	U
129-00-0-----	Pyrene	10	U
85-68-7-----	Butylbenzylphthalate	10	U
91-94-1-----	3,3'-Dichlorobenzidine	10	U
56-55-3-----	Benzo(a)Anthracene	10	U
218-01-9-----	Chrysene	10	U
117-81-7-----	bis(2-Ethylhexyl)Phthalate	1	BJ
117-84-0-----	Di-n-Octyl Phthalate	10	U
205-99-2-----	Benzo(b)Fluoranthene	10	U
207-08-9-----	Benzo(k)Fluoranthene	10	U
50-32-8-----	Benzo(a)Pyrene	10	U
193-39-5-----	Indeno(1,2,3-cd)Pyrene	10	U
53-70-3-----	Dibenz(a,h)Anthracene	10	U
191-24-2-----	Benzo(g,h,i)Perylene	10	U

(1) - Cannot be separated from Diphenylamine

1B
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

ORIGINAL
CKY26
(P-6)

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU

Case No.: 18347

SAS No.: _____

SDG No.: CKY14

Matrix: (soil/water) WATER

Lab Sample ID: 499449

Sample wt/vol: 1000 (g/mL) ML

Lab File ID: GH099449A07

Level: (low/med) LOW

Date Received: 06/24/92

% Moisture: _____ decanted: (Y/N)

Date Extracted: 06/25/92

Concentrated Extract Volume: 1000 (uL)

Date Analyzed: 06/30/92

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) N pH:

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/L

Q

CAS NO.	COMPOUND	10	U
108-95-2-----	Phenol	10	U
111-44-4-----	bis(2-Chloroethyl)Ether	10	U
95-57-8-----	2-Chlorophenol	10	U
541-73-1-----	1,3-Dichlorobenzene	10	U
106-46-7-----	1,4-Dichlorobenzene	10	U
95-50-1-----	1,2-Dichlorobenzene	10	U
95-48-7-----	2-Methylphenol	10	U
108-60-1-----	2,2'-Oxybis(1-Chloropropane)	10	U
106-44-5-----	4-Methylphenol	10	U
621-64-7-----	N-Nitroso-Di-n-Propylamine	10	U
67-72-1-----	Hexachloroethane	10	U
98-95-3-----	Nitrobenzene	10	U
78-59-1-----	Isophorone	10	U
88-75-5-----	2-Nitrophenol	10	U
105-67-9-----	2,4-Dimethylphenol	10	U
111-91-1-----	bis(2-Chloroethoxy)Methane	10	U
120-83-2-----	2,4-Dichlorophenol	10	U
120-82-1-----	1,2,4-Trichlorobenzene	10	U
91-20-3-----	Naphthalene	10	U
106-47-8-----	4-Chloroaniline	10	U
87-68-3-----	Hexachlorobutadiene	10	U
59-50-7-----	4-Chloro-3-Methylphenol	10	U
91-57-6-----	2-Methylnaphthalene	10	U
77-47-4-----	Hexachlorocyclopentadiene	10	U
88-06-2-----	2,4,6-Trichlorophenol	10	U
95-95-4-----	2,4,5-Trichlorophenol	25	U
91-58-7-----	2-Chloronaphthalene	10	U
88-74-4-----	2-Nitroaniline	25	U
131-11-3-----	Dimethyl Phthalate	10	U
208-96-8-----	Acenaphthylene	10	U
606-20-2-----	2,6-Dinitrotoluene	10	U
99-09-2-----	3-Nitroaniline	25	U
83-32-9-----	Acenaphthene	10	U

1C
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CKY26

Lab Name: COMPUCHEM, RTPContract: 68D10083Lab Code: COMPU Case No.: 18347SAS No.: _____ SDG No.: CKY14Matrix: (soil/water) WATERLab Sample ID: 499449Sample wt/vol: 1000 (g/mL) MLLab File ID: GH099449A07Level: (low/med) LOWDate Received: 06/24/92% Moisture: _____ decanted: (Y/N) Date Extracted: 06/25/92Concentrated Extract Volume: 1000 (uL)Date Analyzed: 06/30/92Injection Volume: 2.0(uL)Dilution Factor: 1.0GPC Cleanup: (Y/N) N pH: _____CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/L Q

CAS NO.	COMPOUND	25	U
51-28-5-----	2,4-Dinitrophenol	25	U
100-02-7-----	4-Nitrophenol	25	U
132-64-9-----	Dibenzofuran	10	U
121-14-2-----	2,4-Dinitrotoluene	10	U
84-66-2-----	Diethylphthalate	10	U
7005-72-3-----	4-Chlorophenyl-phenylether	10	U
86-73-7-----	Fluorene	10	U
100-01-6-----	4-Nitroaniline	25	U
534-52-1-----	4,6-Dinitro-2-Methylphenol	25	U
86-30-6-----	N-Nitrosodiphenylamine (1)	10	U
101-55-3-----	4-Bromophenyl-phenylether	10	U
118-74-1-----	Hexachlorobenzene	10	U
87-86-5-----	Pentachlorophenol	25	U
85-01-8-----	Phenanthrene	10	U
120-12-7-----	Anthracene	10	U
86-74-8-----	Carbazole	10	U
84-74-2-----	Di-n-Butylphthalate	10	U
206-44-0-----	Fluoranthene	10	U
129-00-0-----	Pyrene	10	U
85-68-7-----	Butylbenzylphthalate	10	U
91-94-1-----	3,3'-Dichlorobenzidine	10	U
56-55-3-----	Benzo(a)Anthracene	10	U
218-01-9-----	Chrysene	10	U
117-81-7-----	bis(2-Ethylhexyl)Phthalate	2	BJ
117-84-0-----	Di-n-Octyl Phthalate	10	U
205-99-2-----	Benzo(b)Fluoranthene	10	U
207-08-9-----	Benzo(k)Fluoranthene	10	U
50-32-8-----	Benzo(a)Pyrene	10	U
193-39-5-----	Indeno(1,2,3-cd) Pyrene	10	U
53-70-3-----	Dibenz(a,h)Anthracene	10	U
191-24-2-----	Benzo(g,h,i)Perylene	10	U

(1) - Cannot be separated from Diphenylamine

1B
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

(P-60)
FINAL

CKY28

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU

Case No.: 18347

SAS No.: _____

SDG No.: CKY28

Matrix: (soil/water) SOIL

Lab Sample ID: 499452

Sample wt/vol: 30.1 (g/mL) G

Lab File ID: GR099452A57

Level: (low/med) LOW

Date Received: 06/24/92

% Moisture: 19 decanted: (Y/N) N

Date Extracted: 07/07/92

Concentrated Extract Volume: 500.0 (uL)

Date Analyzed: 07/09/92

Injection Volume: 2.0(uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 7.2

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG

Q

CAS NO.	COMPOUND			
108-95-2-----	Phenol	410	U	
111-44-4-----	bis(2-Chloroethyl) Ether	410	U	
95-57-8-----	2-Chlorophenol	410	U	
541-73-1-----	1,3-Dichlorobenzene	410	U	
106-46-7-----	1,4-Dichlorobenzene	410	U	
95-50-1-----	1,2-Dichlorobenzene	410	U	
95-48-7-----	2-Methylphenol	410	U	
108-60-1-----	2,2'-Oxybis(1-Chloropropane)	410	U	
106-44-5-----	4-Methylphenol	410	U	
621-64-7-----	N-Nitroso-Di-n-Propylamine	410	U	
67-72-1-----	Hexachloroethane	410	U	
98-95-3-----	Nitrobenzene	410	U	
78-59-1-----	Isophorone	410	U	
88-75-5-----	2-Nitrophenol	410	U	
105-67-9-----	2,4-Dimethylphenol	410	U	
111-91-1-----	bis(2-Chloroethoxy) Methane	410	U	
120-83-2-----	2,4-Dichlorophenol	410	U	
120-82-1-----	1,2,4-Trichlorobenzene	410	U	
91-20-3-----	Naphthalene	410	U	
106-47-8-----	4-Chloroaniline	410	U	
87-68-3-----	Hexachlorobutadiene	410	U	
59-50-7-----	4-Chloro-3-Methylphenol	410	U	
91-57-6-----	2-Methylnaphthalene	410	U	
77-47-4-----	Hexachlorocyclopentadiene	410	U	
88-06-2-----	2,4,6-Trichlorophenol	410	U	
95-95-4-----	2,4,5-Trichlorophenol	980	U	
91-58-7-----	2-Chloronaphthalene	410	U	
88-74-4-----	2-Nitroaniline	980	U	
131-11-3-----	Dimethyl Phthalate	410	U	
208-96-8-----	Acenaphthylene	410	U	
606-20-2-----	2,6-Dinitrotoluene	410	U	
99-09-2-----	3-Nitroaniline	980	U	
83-32-9-----	Acenaphthene	410	U	

1C
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

*(Signature)
(Red)*

CKY28

Lab Name: COMPUCHEM RTP Contract: 68D10083

Lab Code: COMPU Case No.: 18347 SAS No.: _____ SDG No.: CKY28

Matrix: (soil/water) SOIL Lab Sample ID: 499452

Sample wt/vol: 30.1 (g/mL) G Lab File ID: GR099452A57

Level: (low/med) LOW Date Received: 06/24/92

% Moisture: 19 decanted: (Y/N) N Date Extracted: 07/07/92

Concentrated Extract Volume: 500.0 (uL) Date Analyzed: 07/09/92

Injection Volume: 2.0 (uL) Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 7.2

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG Q

51-28-5-----	2,4-Dinitrophenol	980	U
100-02-7-----	4-Nitrophenol	980	U
132-64-9-----	Dibenzofuran	410	U
121-14-2-----	2,4-Dinitrotoluene	410	U
84-66-2-----	Diethylphthalate	410	U
7005-72-3-----	4-Chlorophenyl-phenylether	410	U
86-73-7-----	Fluorene	410	U
100-01-6-----	4-Nitroaniline	980	U
534-52-1-----	4,6-Dinitro-2-Methylphenol	980	U
86-30-6-----	N-Nitrosodiphenylamine (1)	410	U
101-55-3-----	4-Bromophenyl-phenylether	410	U
118-74-1-----	Hexachlorobenzene	410	U
87-86-5-----	Pentachlorophenol	980	U
85-01-8-----	Phenanthrene	410	U
120-12-7-----	Anthracene	410	U
86-74-8-----	Carbazole	410	U
84-74-2-----	Di-n-Butylphthalate	410	U
206-44-0-----	Fluoranthene	410	U
129-00-0-----	Pyrene	410	U
85-68-7-----	Butylbenzylphthalate	410	U
91-94-1-----	3,3'-Dichlorobenzidine	410	U
56-55-3-----	Benzo(a)Anthracene	410	U
218-01-9-----	Chrysene	410	U
117-81-7-----	bis(2-Ethylhexyl) Phthalate	410	U
117-84-0-----	Di-n-Octyl Phthalate	410	U
205-99-2-----	Benzo(b)Fluoranthene	410	U
207-08-9-----	Benzo(k)Fluoranthene	410	U
50-32-8-----	Benzo(a)Pyrene	410	U
193-39-5-----	Indeno(1,2,3-cd) Pyrene	410	U
53-70-3-----	Dibenz(a,h)Anthracene	410	U
191-24-2-----	Benzo(g,h,i)Perylene	410	U

(1) - Cannot be separated from Diphenylamine

1B
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CKY29

Lab Name: COMPUCHEM,RTP

Contract: 68D10083

Lab Code: COMPU

Case No.: 18347

SAS No.: _____

SDG No.: CKY28

Matrix: (soil/water) SOIL

Lab Sample ID: 499459

Sample wt/vol: 30.7 (g/mL) G

Lab File ID: G2R99459A05

Level: (low/med) LOW

Date Received: 06/24/92

% Moisture: 17 decanted: (Y/N) N

Date Extracted: 07/10/92

Concentrated Extract Volume: 500.0 (uL)

Date Analyzed: 07/15/92

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 6.6

CONCENTRATION UNITS:

CAS NO.	COMPOUND	(ug/L or ug/Kg) UG/KG	Q
---------	----------	-----------------------	---

108-95-2-----	Phenol	390	U
111-44-4-----	bis(2-Chloroethyl)Ether	390	U
95-57-8-----	2-Chlorophenol	390	U
541-73-1-----	1,3-Dichlorobenzene	390	U
106-46-7-----	1,4-Dichlorobenzene	390	U
95-50-1-----	1,2-Dichlorobenzene	390	U
95-48-7-----	2-Methylphenol	390	U
108-60-1-----	2,2'-Oxybis(1-Chloropropane)	390	U
106-44-5-----	4-Methylphenol	390	U
621-64-7-----	N-Nitroso-Di-n-Propylamine	390	U
67-72-1-----	Hexachloroethane	390	U
98-95-3-----	Nitrobenzene	390	U
78-59-1-----	Isophorone	390	U
88-75-5-----	2-Nitrophenol	390	U
105-67-9-----	2,4-Dimethylphenol	390	U
111-91-1-----	bis(2-Chloroethoxy)Methane	390	U
120-83-2-----	2,4-Dichlorophenol	390	U
120-82-1-----	1,2,4-Trichlorobenzene	390	U
91-20-3-----	Naphthalene	390	U
106-47-8-----	4-Chloroaniline	390	U
87-68-3-----	Hexachlorobutadiene	390	U
59-50-7-----	4-Chloro-3-Methylphenol	390	U
91-57-6-----	2-Methylnaphthalene	92	J
77-47-4-----	Hexachlorocyclopentadiene	390	U
88-06-2-----	2,4,6-Trichlorophenol	390	U
95-95-4-----	2,4,5-Trichlorophenol	940	U
91-58-7-----	2-Chloronaphthalene	390	U
88-74-4-----	2-Nitroaniline	940	U
131-11-3-----	Dimethyl Phthalate	390	U
208-96-8-----	Acenaphthylene	390	U
606-20-2-----	2,6-Dinitrotoluene	390	U
99-09-2-----	3-Nitroaniline	940	U
83-32-9-----	Acenaphthene	390	U

1C
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

ORIG/1991

(Rev)

CKY29

Lab Name: COMPUCHEM, RTPContract: 68D10083Lab Code: COMPU Case No.: 18347SAS No.: _____ SDG No.: CKY28Matrix: (soil/water) SOILLab Sample ID: 499459Sample wt/vol: 30.7 (g/mL) GLab File ID: G2R99459A05Level: (low/med) LOWDate Received: 06/24/92Moisture: 17 decanted: (Y/N) NDate Extracted: 07/10/92Concentrated Extract Volume: 500.0 (uL)Date Analyzed: 07/15/92Injection Volume: 2.0 (uL)Dilution Factor: 1.0GPC Cleanup: (Y/N) Y pH: 6.6

CONCENTRATION UNITS:

(ug/L or ug/Kg) UG/KGQ

CAS NO.	COMPOUND	UG/KG	Q
51-28-5-----	2,4-Dinitrophenol	940	U
100-02-7-----	4-Nitrophenol	940	U
132-64-9-----	Dibenzofuran	390	U
121-14-2-----	2,4-Dinitrotoluene	390	U
84-66-2-----	Diethylphthalate	390	U
7005-72-3-----	4-Chlorophenyl-phenylether	390	U
86-73-7-----	Fluorene	390	U
100-01-6-----	4-Nitroaniline	940	U
534-52-1-----	4,6-Dinitro-2-Methylphenol	940	U
86-30-6-----	N-Nitrosodiphenylamine (1)	390	U
101-55-3-----	4-Bromophenyl-phenylether	390	U
118-74-1-----	Hexachlorobenzene	390	U
87-86-5-----	Pentachlorophenol	940	U
85-01-8-----	Phenanthrene	130	J
120-12-7-----	Anthracene	390	U
86-74-8-----	Carbazole	390	U
84-74-2-----	Di-n-Butylphthalate	390	U
206-44-0-----	Fluoranthene	120	J
129-00-0-----	Pyrene	85	J
85-68-7-----	Butylbenzylphthalate	46	BJ
91-94-1-----	3,3'-Dichlorobenzidine	390	U
56-55-3-----	Benzo(a)Anthracene	110	J
218-01-9-----	Chrysene	120	J
117-81-7-----	bis(2-Ethylhexyl) Phthalate	93	J
117-84-0-----	Di-n-Octyl Phthalate	390	U
205-99-2-----	Benzo(b)Fluoranthene	250	JX
207-08-9-----	Benzo(k)Fluoranthene	250	JX
50-32-8-----	Benzo(a)Pyrene	120	J
193-39-5-----	Indeno(1,2,3-cd) Pyrene	64	J
53-70-3-----	Dibenz(a,h)Anthracene	390	U
191-24-2-----	Benzo(g,h,i)Perylene	390	U

(1) - Cannot be separated from Diphenylamine

1B
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

ORIGINAL
(P&G)

CKY30

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 18347

SAS No.: _____ SDG No.: CKY28

Matrix: (soil/water) SOIL

Lab Sample ID: 499463

Sample wt/vol: 30.0 (g/mL) G

Lab File ID: GH099463A15

Level: (low/med) LOW

Date Received: 06/24/92

% Moisture: 40 decanted: (Y/N) N

Date Extracted: 06/26/92

Concentrated Extract Volume: 500.0 (uL)

Date Analyzed: 07/01/92

Injection Volume: 2.0(uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 5.5

CONCENTRATION UNITS:

(ug/L or ug/Kg) UG/KG

Q

CAS NO.	COMPOUND	Q
108-95-2-----	Phenol	550 U
111-44-4-----	bis(2-Chloroethyl)Ether	550 U
95-57-8-----	2-Chlorophenol	550 U
541-73-1-----	1,3-Dichlorobenzene	550 U
106-46-7-----	1,4-Dichlorobenzene	550 U
95-50-1-----	1,2-Dichlorobenzene	550 U
95-48-7-----	2-Methylphenol	550 U
108-60-1-----	2,2'-Oxybis(1-Chloropropane)	550 U
106-44-5-----	4-Methylphenol	62 J
621-64-7-----	N-Nitroso-Di-n-Propylamine	550 U
67-72-1-----	Hexachloroethane	550 U
98-95-3-----	Nitrobenzene	550 U
78-59-1-----	Isophorone	550 U
88-75-5-----	2-Nitrophenol	550 U
105-67-9-----	2,4-Dimethylphenol	550 U
111-91-1-----	bis(2-Chloroethoxy)Methane	550 U
120-83-2-----	2,4-Dichlorophenol	550 U
120-82-1-----	1,2,4-Trichlorobenzene	550 U
91-20-3-----	Naphthalene	550 U
106-47-8-----	4-Chloroaniline	550 U
87-68-3-----	Hexachlorobutadiene	550 U
59-50-7-----	4-Chloro-3-Methylphenol	550 U
91-57-6-----	2-Methylnaphthalene	550 U
77-47-4-----	Hexachlorocyclopentadiene	550 U
88-06-2-----	2,4,6-Trichlorophenol	550 U
95-95-4-----	2,4,5-Trichlorophenol	1300 U
91-58-7-----	2-Chloronaphthalene	550 U
88-74-4-----	2-Nitroaniline	1300 U
131-11-3-----	Dimethyl Phthalate	550 U
208-96-8-----	Acenaphthylene	550 U
606-20-2-----	2,6-Dinitrotoluene	550 U
99-09-2-----	3-Nitroaniline	1300 U
83-32-9-----	Acenaphthene	550 U

1C
SEMICVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

CKY28
(Red)

Lab Code: COMPU

Case No.: 18347

SAS No.: _____

SDG No.: CKY28

Matrix: (soil/water) SOIL

Lab Sample ID: 499463

Sample wt/vol: 30.0 (g/mL) G

Lab File ID: GH099463A15

Level: (low/med) LOW

Date Received: 06/24/92

% Moisture: 40 decanted: (Y/N) N

Date Extracted: 06/26/92

Concentrated Extract Volume: 500.0 (uL)

Date Analyzed: 07/01/92

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 5.5

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG Q

CAS NO.	COMPOUND	1300	U
51-28-5-----	2,4-Dinitrophenol	1300	U
100-02-7-----	4-Nitrophenol	1300	U
132-64-9-----	Dibenzofuran	550	U
121-14-2-----	2,4-Dinitrotoluene	550	U
84-66-2-----	Diethylphthalate	550	U
7005-72-3-----	4-Chlorophenyl-phenylether	550	U
86-73-7-----	Fluorene	550	U
100-01-6-----	4-Nitroaniline	1300	U
534-52-1-----	4,6-Dinitro-2-Methylphenol	1300	U
86-30-6-----	N-Nitrosodiphenylamine (1)	550	U
101-55-3-----	4-Bromophenyl-phenylether	550	U
118-74-1-----	Hexachlorobenzene	550	U
87-86-5-----	Pentachlorophenol	1300	U
85-01-8-----	Phenanthrene	56	J
120-12-7-----	Anthracene	550	U
86-74-8-----	Carbazole	550	U
84-74-2-----	Di-n-Butylphthalate	550	U
206-44-0-----	Fluoranthene	550	U
129-00-0-----	Pyrene	550	U
85-68-7-----	Butylbenzylphthalate	240	BJ
91-94-1-----	3,3'-Dichlorobenzidine	550	U
56-55-3-----	Benzo(a)Anthracene	550	U
218-01-9-----	Chrysene	550	U
117-81-7-----	bis(2-Ethylhexyl)Phthalate	550	U
117-84-0-----	Di-n-Octyl Phthalate	550	U
205-99-2-----	Benzo(b)Fluoranthene	550	U
207-08-9-----	Benzo(k)Fluoranthene	550	U
50-32-8-----	Benzo(a)Pyrene	550	U
193-39-5-----	Indeno(1,2,3-cd)Pyrene	550	U
53-70-3-----	Dibenz(a,h)Anthracene	550	U
191-24-2-----	Benzo(g,h,i)Perylene	550	U

(1) - Cannot be separated from Diphenylamine

1B
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

OB
FINAL
(Rev)

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 18347

SAS No.: _____ SDG No.: CKY28

Matrix: (soil/water) SOIL

Lab Sample ID: 499464

Sample wt/vol: 30.0 (g/mL) G

Lab File ID: GR099464A57

Level: (low/med) LOW

Date Received: 06/24/92

% Moisture: 23 decanted: (Y/N) N

Date Extracted: 07/07/92

Concentrated Extract Volume: 500.0 (uL)

Date Analyzed: 07/09/92

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 7.1

CONCENTRATION UNITS:

CAS NO.	COMPOUND	(ug/L or ug/Kg) <u>UG/KG</u>	Q
---------	----------	------------------------------	---

108-95-2-----	Phenol	430	U
111-44-4-----	bis(2-Chloroethyl)Ether	430	U
95-57-8-----	2-Chlorophenol	430	U
541-73-1-----	1,3-Dichlorobenzene	430	U
106-46-7-----	1,4-Dichlorobenzene	430	U
95-50-1-----	1,2-Dichlorobenzene	430	U
95-48-7-----	2-Methylphenol	430	U
108-60-1-----	2,2'-Oxybis(1-Chloropropane)	430	U
106-44-5-----	4-Methylphenol	430	U
621-64-7-----	N-Nitroso-Di-n-Propylamine	430	U
67-72-1-----	Hexachloroethane	430	U
98-95-3-----	Nitrobenzene	430	U
78-59-1-----	Isophorone	430	U
88-75-5-----	2-Nitrophenol	430	U
105-67-9-----	2,4-Dimethylphenol	430	U
111-91-1-----	bis(2-Chloroethoxy)Methane	430	U
120-83-2-----	2,4-Dichlorophenol	430	U
120-82-1-----	1,2,4-Trichlorobenzene	430	U
91-20-3-----	Naphthalene	430	U
106-47-8-----	4-Chloroaniline	430	U
87-68-3-----	Hexachlorobutadiene	430	U
59-50-7-----	4-Chloro-3-Methylphenol	430	U
91-57-6-----	2-Methylnaphthalene	430	U
77-47-4-----	Hexachlorocyclopentadiene	430	U
88-06-2-----	2,4,6-Trichlorophenol	430	U
95-95-4-----	2,4,5-Trichlorophenol	1000	U
91-58-7-----	2-Chloronaphthalene	430	U
88-74-4-----	2-Nitroaniline	1000	U
131-11-3-----	Dimethyl Phthalate	430	U
208-96-8-----	Acenaphthylene	430	U
606-20-2-----	2,6-Dinitrotoluene	430	U
99-09-2-----	3-Nitroaniline	1000	U
83-32-9-----	Acenaphthene	430	U

1C
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

ORIGINAL
(med)

CKY31

Lab Name: COMPUCHEM, RTPContract: 68D10083Lab Code: COMPU Case No.: 18347SAS No.: _____ SDG No.: CKY28Matrix: (soil/water) SOILLab Sample ID: 499464Sample wt/vol: 30.0 (g/mL) GLab File ID: GR099464A57Level: (low/med) LOWDate Received: 06/24/92% Moisture: 23 decanted: (Y/N) NDate Extracted: 07/07/92Concentrated Extract Volume: 500.0 (uL)Date Analyzed: 07/09/92Injection Volume: 2.0 (uL)Dilution Factor: 1.0GPC Cleanup: (Y/N) Y pH: 7.1CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG

Q

CAS NO.	COMPOUND		
51-28-5-----	2,4-Dinitrophenol	1000	U
100-02-7-----	4-Nitrophenol	1000	U
132-64-9-----	Dibenzofuran	430	U
121-14-2-----	2,4-Dinitrotoluene	430	U
84-66-2-----	Diethylphthalate	430	U
7005-72-3-----	4-Chlorophenyl-phenylether	430	U
86-73-7-----	Fluorene	430	U
100-01-6-----	4-Nitroaniline	1000	U
534-52-1-----	4,6-Dinitro-2-Methylphenol	1000	U
86-30-6-----	N-Nitrosodiphenylamine (1)	430	U
101-55-3-----	4-Bromophenyl-phenylether	430	U
118-74-1-----	Hexachlorobenzene	430	U
87-86-5-----	Pentachlorophenol	1000	U
85-01-8-----	Phenanthrene	430	U
120-12-7-----	Anthracene	430	U
86-74-8-----	Carbazole	430	U
84-74-2-----	Di-n-Butylphthalate	430	U
206-44-0-----	Fluoranthene	430	U
129-00-0-----	Pyrene	430	U
85-68-7-----	Butylbenzylphthalate	430	U
91-94-1-----	3,3'-Dichlorobenzidine	430	U
56-55-3-----	Benzo(a)Anthracene	430	U
218-01-9-----	Chrysene	430	U
117-81-7-----	bis(2-Ethylhexyl)Phthalate	430	U
117-84-0-----	Di-n-Octyl Phthalate	430	U
205-99-2-----	Benzo(b)Fluoranthene	430	U
207-08-9-----	Benzo(k)Fluoranthene	430	U
50-32-8-----	Benzo(a)Pyrene	430	U
193-39-5-----	Indeno(1,2,3-cd) Pyrene	430	U
53-70-3-----	Dibenz(a,h)Anthracene	430	U
191-24-2-----	Benzo(g,h,i)Perylene	430	U

(1) - Cannot be separated from Diphenylamine

SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

CKY32

Lab Name: COMPUCHEM, RTP Contract: 68D10083Lab Code: COMPU Case No.: 18347 SAS No.: _____ SDG No.: CKY28Matrix: (soil/water) SOILLab Sample ID: 499465Sample wt/vol: 30.0 (g/mL) GLab File ID: GH099465A15Level: (low/med) LOWDate Received: 06/24/92% Moisture: 18 decanted: (Y/N) NDate Extracted: 06/26/92Concentrated Extract Volume: 500.0 (uL)Date Analyzed: 07/01/92Injection Volume: 2.0 (uL)Dilution Factor: 1.0GPC Cleanup: (Y/N) Y pH: 6.5CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG Q

CAS NO. COMPOUND

108-95-2-----	Phenol	400	U
111-44-4-----	bis(2-Chloroethyl) Ether	400	U
95-57-8-----	2-Chlorophenol	400	U
541-73-1-----	1,3-Dichlorobenzene	400	U
106-46-7-----	1,4-Dichlorobenzene	400	U
95-50-1-----	1,2-Dichlorobenzene	400	U
95-48-7-----	2-Methylphenol	400	U
108-60-1-----	2,2'-Oxybis(1-Chloropropane)	400	U
106-44-5-----	4-Methylphenol	400	U
621-64-7-----	N-Nitroso-Di-n-Propylamine	400	U
67-72-1-----	Hexachloroethane	400	U
98-95-3-----	Nitrobenzene	400	U
78-59-1-----	Isophorone	400	U
88-75-5-----	2-Nitrophenol	400	U
105-67-9-----	2,4-Dimethylphenol	400	U
111-91-1-----	bis(2-Chloroethoxy)Methane	400	U
120-83-2-----	2,4-Dichlorophenol	400	U
120-82-1-----	1,2,4-Trichlorobenzene	400	U
91-20-3-----	Naphthalene	400	U
106-47-8-----	4-Chloroaniline	400	U
87-68-3-----	Hexachlorobutadiene	400	U
59-50-7-----	4-Chloro-3-Methylphenol	400	U
91-57-6-----	2-Methylnaphthalene	400	U
77-47-4-----	Hexachlorocyclopentadiene	400	U
88-06-2-----	2,4,6-Trichlorophenol	400	U
95-95-4-----	2,4,5-Trichlorophenol	980	U
91-58-7-----	2-Chloronaphthalene	400	U
88-74-4-----	2-Nitroaniline	980	U
131-11-3-----	Dimethyl Phthalate	400	U
208-96-8-----	Acenaphthylene	400	U
606-20-2-----	2,6-Dinitrotoluene	400	U
99-09-2-----	3-Nitroaniline	980	U
83-32-9-----	Acenaphthene	400	U

1C
SEMICVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CKY32

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 18347

SAS No.: _____ SDG No.: CKY28

Matrix: (soil/water) SOIL

Lab Sample ID: 499465

Sample wt/vol: 30.0 (g/mL) G

Lab File ID: GH099465A15

Level: (low/med) LOW

Date Received: 06/24/92

% Moisture: 18 decanted: (Y/N) N

Date Extracted: 06/26/92

Concentrated Extract Volume: 500.0 (uL)

Date Analyzed: 07/01/92

Injection Volume: 2.0(uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 6.5

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG Q

CAS NO.	COMPOUND	Q
51-28-5-----	2,4-Dinitrophenol	980 U
100-02-7-----	4-Nitrophenol	980 U
132-64-9-----	Dibenzofuran	400 U
121-14-2-----	2,4-Dinitrotoluene	400 U
84-66-2-----	Diethylphthalate	400 U
7005-72-3-----	4-Chlorophenyl-phenylether	400 U
86-73-7-----	Fluorene	400 U
100-01-6-----	4-Nitroaniline	980 U
534-52-1-----	4,6-Dinitro-2-Methylphenol	980 U
86-30-6-----	N-Nitrosodiphenylamine (1)	400 U
101-55-3-----	4-Bromophenyl-phenylether	400 U
118-74-1-----	Hexachlorobenzene	400 U
87-86-5-----	Pentachlorophenol	980 U
85-01-8-----	Phenanthrene	400 U
120-12-7-----	Anthracene	400 U
86-74-8-----	Carbazole	400 U
84-74-2-----	Di-n-Butylphthalate	400 U
206-44-0-----	Fluoranthene	400 U
129-00-0-----	Pyrene	400 U
85-68-7-----	Butylbenzylphthalate	180 BJ
91-94-1-----	3,3'-Dichlorobenzidine	400 U
56-55-3-----	Benzo(a)Anthracene	400 U
218-01-9-----	Chrysene	400 U
117-81-7-----	bis(2-Ethylhexyl) Phthalate	400 U
117-84-0-----	Di-n-Octyl Phthalate	400 U
205-99-2-----	Benzo(b)Fluoranthene	400 U
207-08-9-----	Benzo(k)Fluoranthene	400 U
50-32-8-----	Benzo(a)Pyrene	400 U
193-39-5-----	Indeno(1,2,3-cd)Pyrene	400 U
53-70-3-----	Dibenz(a,h)Anthracene	400 U
191-24-2-----	Benzo(g,h,i)Perylene	400 U

(1) - Cannot be separated from Diphenylamine

1B
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

Lab Name: COMPUCHEM.RTPContract: 68D10083CKY35
^{ORIGINAL}
(Red)Lab Code: COMPU Case No.: 18347SAS No.: _____ SDG No.: CKY28Matrix: (soil/water) SOILLab Sample ID: 499466Sample wt/vol: 30.0 (g/mL) GLab File ID: GH099466A15Level: (low/med) LOWDate Received: 06/24/92% Moisture: 11 decanted: (Y/N) NDate Extracted: 06/26/92Concentrated Extract Volume: 500.0 (uL)Date Analyzed: 07/02/92Injection Volume: 2.0 (uL)Dilution Factor: 1.0GPC Cleanup: (Y/N) Y pH: 5.9

CONCENTRATION UNITS:

CAS NO. COMPOUND (ug/L or ug/Kg) UG/KG Q

108-95-2-----	Phenol	370	U
111-44-4-----	bis(2-Chloroethyl) Ether	370	U
95-57-8-----	2-Chlorophenol	370	U
541-73-1-----	1,3-Dichlorobenzene	370	U
106-46-7-----	1,4-Dichlorobenzene	370	U
95-50-1-----	1,2-Dichlorobenzene	370	U
95-48-7-----	2-Methylphenol	370	U
108-60-1-----	2,2'-Oxybis(1-Chloropropane)	370	U
106-44-5-----	4-Methylphenol	370	U
621-64-7-----	N-Nitroso-Di-n-Propylamine	370	U
67-72-1-----	Hexachloroethane	370	U
98-95-3-----	Nitrobenzene	370	U
78-59-1-----	Isophorone	370	U
88-75-5-----	2-Nitrophenol	370	U
105-67-9-----	2,4-Dimethylphenol	370	U
111-91-1-----	bis(2-Chloroethoxy) Methane	370	U
120-83-2-----	2,4-Dichlorophenol	370	U
120-82-1-----	1,2,4-Trichlorobenzene	370	U
91-20-3-----	Naphthalene	370	U
106-47-8-----	4-Chloroaniline	370	U
87-68-3-----	Hexachlorobutadiene	370	U
59-50-7-----	4-Chloro-3-Methylphenol	370	U
91-57-6-----	2-Methylnaphthalene	370	U
77-47-4-----	Hexachlorocyclopentadiene	370	U
88-06-2-----	2,4,6-Trichlorophenol	370	U
95-95-4-----	2,4,5-Trichlorophenol	900	U
91-58-7-----	2-Chloronaphthalene	370	U
88-74-4-----	2-Nitroaniline	900	U
131-11-3-----	Dimethyl Phthalate	370	U
208-96-8-----	Acenaphthylene	370	U
606-20-2-----	2,6-Dinitrotoluene	370	U
99-09-2-----	3-Nitroaniline	900	U
83-32-9-----	Acenaphthene	370	U

FORM I SV-1

3/90

1C
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CKY33 (Rej)

Lab Name: COMPUCHEM, RTPContract: 68D10083Lab Code: COMPUCase No.: 18347

SAS No.: _____

SDG No.: CKY28Matrix: (soil/water) SOILLab Sample ID: 499466Sample wt/vol: 30.0 (g/mL) GLab File ID: GH099466A15Level: (low/med) LOWDate Received: 06/24/92% Moisture: 11 decanted: (Y/N) NDate Extracted: 06/26/92Concentrated Extract Volume: 500.0 (uL)Date Analyzed: 07/02/92Injection Volume: 2.0(uL)Dilution Factor: 1.0GPC Cleanup: (Y/N) Y pH: 5.9CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG Q

CAS NO.	COMPOUND	Q
51-28-5-----	2,4-Dinitrophenol	900 U
100-02-7-----	4-Nitrophenol	900 U
132-64-9-----	Dibenzofuran	370 U
121-14-2-----	2,4-Dinitrotoluene	370 U
84-66-2-----	Diethylphthalate	370 U
7005-72-3-----	4-Chlorophenyl-phenylether	370 U
86-73-7-----	Fluorene	370 U
100-01-6-----	4-Nitroaniline	900 U
534-52-1-----	4,6-Dinitro-2-Methylphenol	900 U
86-30-6-----	N-Nitrosodiphenylamine (1)	370 U
101-55-3-----	4-Bromophenyl-phenylether	370 U
118-74-1-----	Hexachlorobenzene	370 U
87-86-5-----	Pentachlorophenol	900 U
85-01-8-----	Phenanthrene	370 U
120-12-7-----	Anthracene	370 U
86-74-8-----	Carbazole	370 U
84-74-2-----	Di-n-Butylphthalate	370 U
206-44-0-----	Fluoranthene	370 U
129-00-0-----	Pyrene	370 U
85-68-7-----	Butylbenzylphthalate	160 BJ
91-94-1-----	3,3'-Dichlorobenzidine	370 U
56-55-3-----	Benzo(a)Anthracene	370 U
218-01-9-----	Chrysene	370 U
117-81-7-----	bis(2-Ethylhexyl) Phthalate	38 J
117-84-0-----	Di-n-Octyl Phthalate	370 U
205-99-2-----	Benzo(b) Fluoranthene	370 U
207-08-9-----	Benzo(k) Fluoranthene	370 U
50-32-8-----	Benzo(a) Pyrene	370 U
193-39-5-----	Indeno(1,2,3-cd) Pyrene	370 U
53-70-3-----	Dibenz(a,h) Anthracene	370 U
191-24-2-----	Benzo(g,h,i) Perylene	370 U

(1) - Cannot be separated from Diphenylamine

1B
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

(R-91)

CKY34

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 18347

SAS No.: _____ SDG No.: CKY28

Matrix: (soil/water) SOIL

Lab Sample ID: 499467

Sample wt/vol: 30.0 (g/mL) G

Lab File ID: GR099467B57

Level: (low/med) LOW

Date Received: 06/24/92

% Moisture: 13 decanted: (Y/N) N

Date Extracted: 07/07/92

Concentrated Extract Volume: 500.0 (uL)

Date Analyzed: 07/09/92

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 6.5

CONCENTRATION UNITS:

(ug/L or ug/Kg) UG/KG

Q

CAS NO.	COMPOUND	380	U
108-95-2-----	Phenol	380	U
111-44-4-----	bis(2-Chloroethyl)Ether	380	U
95-57-8-----	2-Chlorophenol	380	U
541-73-1-----	1,3-Dichlorobenzene	380	U
106-46-7-----	1,4-Dichlorobenzene	380	U
95-50-1-----	1,2-Dichlorobenzene	380	U
95-48-7-----	2-Methylphenol	380	U
108-60-1-----	2,2'-Oxybis(1-Chloropropane)	380	U
106-44-5-----	4-Methylphenol	380	U
621-64-7-----	N-Nitroso-Di-n-Propylamine	380	U
67-72-1-----	Hexachloroethane	380	U
98-95-3-----	Nitrobenzene	380	U
78-59-1-----	Isophorone	380	U
88-75-5-----	2-Nitrophenol	380	U
105-67-9-----	2,4-Dimethylphenol	380	U
111-91-1-----	bis(2-Chloroethoxy)Methane	380	U
120-83-2-----	2,4-Dichlorophenol	380	U
120-82-1-----	1,2,4-Trichlorobenzene	380	U
91-20-3-----	Naphthalene	380	U
106-47-8-----	4-Chloroaniline	380	U
87-68-3-----	Hexachlorobutadiene	380	U
59-50-7-----	4-Chloro-3-Methylphenol	380	U
91-57-6-----	2-Methylnaphthalene	380	U
77-47-4-----	Hexachlorocyclopentadiene	380	U
88-06-2-----	2,4,6-Trichlorophenol	380	U
95-95-4-----	2,4,5-Trichlorophenol	920	U
91-58-7-----	2-Chloronaphthalene	380	U
88-74-4-----	2-Nitroaniline	920	U
131-11-3-----	Dimethyl Phthalate	380	U
208-96-8-----	Acenaphthylene	380	U
606-20-2-----	2,6-Dinitrotoluene	380	U
99-09-2-----	3-Nitroaniline	920	U
83-32-9-----	Acenaphthene	380	U

1C
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

ORIGINAL
(red)

CKY34

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 18347

SAS No.: _____ SDG No.: CKY28

Matrix: (soil/water) SOIL

Lab Sample ID: 499467

Sample wt/vol: 30.0 (g/mL) G

Lab File ID: GR099467B57

Level: (low/med) LOW

Date Received: 06/24/92

% Moisture: 13 decanted: (Y/N) N

Date Extracted: 07/07/92

Concentrated Extract Volume: 500.0 (uL)

Date Analyzed: 07/09/92

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 6.5

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG Q

CAS NO.	COMPOUND			
51-28-5-----	2,4-Dinitrophenol	920	U	
100-02-7-----	4-Nitrophenol	920	U	
132-64-9-----	Dibenzofuran	380	U	
121-14-2-----	2,4-Dinitrotoluene	380	U	
84-66-2-----	Diethylphthalate	380	U	
7005-72-3-----	4-Chlorophenyl-phenylether	380	U	
86-73-7-----	Fluorene	380	U	
100-01-6-----	4-Nitroaniline	920	U	
534-52-1-----	4,6-Dinitro-2-Methylphenol	920	U	
86-30-6-----	N-Nitrosodiphenylamine (1)	380	U	
101-55-3-----	4-Bromophenyl-phenylether	380	U	
118-74-1-----	Hexachlorobenzene	380	U	
87-86-5-----	Pentachlorophenol	920	U	
85-01-8-----	Phenanthrene	380	U	
120-12-7-----	Anthracene	380	U	
86-74-8-----	Carbazole	380	U	
84-74-2-----	Di-n-Butylphthalate	380	U	
206-44-0-----	Fluoranthene	39	J	
129-00-0-----	Pyrene	380	U	
85-68-7-----	Butylbenzylphthalate	380	U	
91-94-1-----	3,3'-Dichlorobenzidine	380	U	
56-55-3-----	Benzo(a)Anthracene	380	U	
218-01-9-----	Chrysene	380	U	
117-81-7-----	bis(2-Ethylhexyl)Phthalate	380	U	
117-84-0-----	Di-n-Octyl Phthalate	380	U	
205-99-2-----	Benzo(b)Fluoranthene	380	U	
207-08-9-----	Benzo(k)Fluoranthene	380	U	
50-32-8-----	Benzo(a)Pyrene	380	U	
193-39-5-----	Indeno(1,2,3-cd)Pyrene	380	U	
53-70-3-----	Dibenz(a,h)Anthracene	380	U	
191-24-2-----	Benzo(g,h,i)Perylene	380	U	

(1) - Cannot be separated from Diphenylamine

1B
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

ORIGINAL
(Rev)

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

CKY35

Lab Code: COMPU

Case No.: 18347

SAS No.: _____

SDG No.: CKY28

Matrix: (soil/water) SOIL

Lab Sample ID: 499468

Sample wt/vol: 30.0 (g/mL) G

Lab File ID: GR099468A57

Level: (low/med) LOW

Date Received: 06/24/92

% Moisture: 3 decanted: (Y/N) N

Date Extracted: 07/07/92

Concentrated Extract Volume: 500.0 (uL)

Date Analyzed: 07/09/92

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 6.9

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG **Q**

CAS NO.	COMPOUND			
108-95-2-----	Phenol	340	U	
111-44-4-----	bis(2-Chloroethyl) Ether	340	U	
95-57-8-----	2-Chlorophenol	340	U	
541-73-1-----	1,3-Dichlorobenzene	340	U	
106-46-7-----	1,4-Dichlorobenzene	340	U	
95-50-1-----	1,2-Dichlorobenzene	340	U	
95-48-7-----	2-Methylphenol	340	U	
108-60-1-----	2,2'-Oxybis(1-Chloropropane)	340	U	
106-44-5-----	4-Methylphenol	340	U	
621-64-7-----	N-Nitroso-Di-n-Propylamine	340	U	
67-72-1-----	Hexachloroethane	340	U	
98-95-3-----	Nitrobenzene	340	U	
78-59-1-----	Isophorone	340	U	
88-75-5-----	2-Nitrophenol	340	U	
105-67-9-----	2,4-Dimethylphenol	340	U	
111-91-1-----	bis(2-Chloroethoxy) Methane	340	U	
120-83-2-----	2,4-Dichlorophenol	340	U	
120-82-1-----	1,2,4-Trichlorobenzene	340	U	
91-20-3-----	Naphthalene	340	U	
106-47-8-----	4-Chloroaniline	340	U	
87-68-3-----	Hexachlorobutadiene	340	U	
59-50-7-----	4-Chloro-3-Methylphenol	340	U	
91-57-6-----	2-Methylnaphthalene	340	U	
77-47-4-----	Hexachlorocyclopentadiene	340	U	
88-06-2-----	2,4,6-Trichlorophenol	340	U	
95-95-4-----	2,4,5-Trichlorophenol	820	U	
91-58-7-----	2-Chloronaphthalene	340	U	
88-74-4-----	2-Nitroaniline	820	U	
131-11-3-----	Dimethyl Phthalate	340	U	
208-96-8-----	Acenaphthylene	340	U	
606-20-2-----	2,6-Dinitrotoluene	340	U	
99-09-2-----	3-Nitroaniline	820	U	
83-32-9-----	Acenaphthene	340	U	

1C
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CKY35

Lab Name: COMPUCHEM RTPContract: 68D10083Lab Code: COMPU Case No.: 18347SAS No.: _____ SDG No.: CKY28Matrix: (soil/water) SOILLab Sample ID: 499468Sample wt/vol: 30.0 (g/mL) GLab File ID: GR099468A57Level: (low/med) LOWDate Received: 06/24/92Moisture: 3 decanted: (Y/N) NDate Extracted: 07/07/92Concentrated Extract Volume: 500.0 (uL)Date Analyzed: 07/09/92Injection Volume: 2.0 (uL)Dilution Factor: 1.0GPC Cleanup: (Y/N) Y pH: 6.9

CONCENTRATION UNITS:

(ug/L or ug/Kg) UG/KG

Q

CAS NO.	COMPOUND	820	U
51-28-5-----	2,4-Dinitrophenol	820	U
100-02-7-----	4-Nitrophenol	820	U
132-64-9-----	Dibenzofuran	340	U
121-14-2-----	2,4-Dinitrotoluene	340	U
84-66-2-----	Diethylphthalate	340	U
7005-72-3-----	4-Chlorophenyl-phenylether	340	U
86-73-7-----	Fluorene	340	U
100-01-6-----	4-Nitroaniline	820	U
534-52-1-----	4,6-Dinitro-2-Methylphenol	820	U
86-30-6-----	N-Nitrosodiphenylamine (1)	340	U
101-55-3-----	4-Bromophenyl-phenylether	340	U
118-74-1-----	Hexachlorobenzene	340	U
87-86-5-----	Pentachlorophenol	820	U
85-01-8-----	Phenanthrene	47	J
120-12-7-----	Anthracene	340	U
86-74-8-----	Carbazole	340	U
84-74-2-----	Di-n-Butylphthalate	340	U
206-44-0-----	Fluoranthene	340	U
129-00-0-----	Pyrene	340	U
85-68-7-----	Butylbenzylphthalate	340	U
91-94-1-----	3,3'-Dichlorobenzidine	340	U
56-55-3-----	Benzo(a)Anthracene	340	U
218-01-9-----	Chrysene	340	U
117-81-7-----	bis(2-Ethylhexyl)Phthalate	340	U
117-84-0-----	Di-n-Octyl Phthalate	340	U
205-99-2-----	Benzo(b)Fluoranthene	340	U
207-08-9-----	Benzo(k)Fluoranthene	340	U
50-32-8-----	Benzo(a)Pyrene	340	U
193-39-5-----	Indeno(1,2,3-cd)Pyrene	340	U
53-70-3-----	Dibenz(a,h)Anthracene	340	U
191-24-2-----	Benzo(g,h,i)Perylene	340	U

(1) - Cannot be separated from Diphenylamine

SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

CKY36

Lab Name: COMPUCHEM, RTPContract: 68D10083Lab Code: COMPUCase No.: 18347

SAS No.: _____

SDG No.: CKY28Matrix: (soil/water) SOILLab Sample ID: 499470Sample wt/vol: 30.0 (g/mL) GLab File ID: GR099470A57Level: (low/med) LOWDate Received: 06/24/92% Moisture: 10 decanted: (Y/N) NDate Extracted: 07/07/92Concentrated Extract Volume: 500.0 (uL)Date Analyzed: 07/09/92Injection Volume: 2.0(uL)Dilution Factor: 1.0GPC Cleanup: (Y/N) Y pH: 7.0

CONCENTRATION UNITS:

(ug/L or ug/Kg) UG/KG

Q

CAS NO.	COMPOUND	Q
108-95-2-----	Phenol	370 U
111-44-4-----	bis(2-Chloroethyl)Ether	370 U
95-57-8-----	2-Chlorophenol	370 U
541-73-1-----	1,3-Dichlorobenzene	370 U
106-46-7-----	1,4-Dichlorobenzene	370 U
95-50-1-----	1,2-Dichlorobenzene	370 U
95-48-7-----	2-Methylphenol	370 U
108-60-1-----	2,2'-Oxybis(1-Chloropropane)	370 U
106-44-5-----	4-Methylphenol	370 U
621-64-7-----	N-Nitroso-Di-n-Propylamine	370 U
67-72-1-----	Hexachloroethane	370 U
98-95-3-----	Nitrobenzene	370 U
78-59-1-----	Isophorone	370 U
88-75-5-----	2-Nitrophenol	370 U
105-67-9-----	2,4-Dimethylphenol	370 U
111-91-1-----	bis(2-Chloroethoxy)Methane	370 U
120-83-2-----	2,4-Dichlorophenol	370 U
120-82-1-----	1,2,4-Trichlorobenzene	370 U
91-20-3-----	Naphthalene	370 U
106-47-8-----	4-Chloroaniline	370 U
87-68-3-----	Hexachlorobutadiene	370 U
59-50-7-----	4-Chloro-3-Methylphenol	370 U
91-57-6-----	2-Methylnaphthalene	370 U
77-47-4-----	Hexachlorocyclopentadiene	370 U
88-06-2-----	2,4,6-Trichlorophenol	370 U
95-95-4-----	2,4,5-Trichlorophenol	890 U
91-58-7-----	2-Choronaphthalene	370 U
88-74-4-----	2-Nitroaniline	890 U
131-11-3-----	Dimethyl Phthalate	370 U
208-96-8-----	Acenaphthylene	370 U
606-20-2-----	2,6-Dinitrotoluene	370 U
99-09-2-----	3-Nitroaniline	890 U
83-32-9-----	Acenaphthene	370 U

1C
SEMOVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

(original)
(red)

CKY36

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 18347

SAS No.: _____

SDG No.: CKY28

Matrix: (soil/water) SOIL

Lab Sample ID: 499470

Sample wt/vol: 30.0 (g/mL) G

Lab File ID: GR099470A57

Level: (low/med) LOW

Date Received: 06/24/92

% Moisture: 10 decanted: (Y/N) N

Date Extracted: 07/07/92

Concentrated Extract Volume: 500.0 (uL)

Date Analyzed: 07/09/92

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 7.0

CONCENTRATION UNITS:

(ug/L or ug/Kg) UG/KG

Q

CAS NO.	COMPOUND	890	U
51-28-5-----	2,4-Dinitrophenol	890	U
100-02-7-----	4-Nitrophenol	890	U
132-64-9-----	Dibenzofuran	370	U
121-14-2-----	2,4-Dinitrotoluene	370	U
84-66-2-----	Diethylphthalate	370	U
7005-72-3-----	4-Chlorophenyl-phenylether	370	U
86-73-7-----	Fluorene	370	U
100-01-6-----	4-Nitroaniline	890	U
534-52-1-----	4,6-Dinitro-2-Methylphenol	890	U
86-30-6-----	N-Nitrosodiphenylamine (1)	370	U
101-55-3-----	4-Bromophenyl-phenylether	370	U
118-74-1-----	Hexachlorobenzene	370	U
87-86-5-----	Pentachlorophenol	890	U
85-01-8-----	Phenanthrene	370	U
120-12-7-----	Anthracene	370	U
86-74-8-----	Carbazole	370	U
84-74-2-----	Di-n-Butylphthalate	370	U
206-44-0-----	Fluoranthene	370	U
129-00-0-----	Pyrene	370	U
85-68-7-----	Butylbenzylphthalate	370	U
91-94-1-----	3,3'-Dichlorobenzidine	370	U
56-55-3-----	Benzo(a)Anthracene	370	U
218-01-9-----	Chrysene	370	U
117-81-7-----	bis(2-Ethylhexyl)Phthalate	370	U
117-84-0-----	Di-n-Octyl Phthalate	370	U
205-99-2-----	Benzo(b)Fluoranthene	370	U
207-08-9-----	Benzo(k)Fluoranthene	370	U
50-32-8-----	Benzo(a)Pyrene	370	U
193-39-5-----	Indeno(1,2,3-cd) Pyrene	370	U
53-70-3-----	Dibenz(a,h)Anthracene	370	U
191-24-2-----	Benzo(g,h,i)Perylene	370	U

(1) - Cannot be separated from Diphenylamine

1B
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

Lab Name: COMPUCHEM.RTPContract: 68D10083CKY37Lab Code: COMPU Case No.: 18347 SAS No.: _____ SDG No.: CKY28Matrix: (soil/water) SOIL Lab Sample ID: 499471Sample wt/vol: 30.0 (g/mL) G Lab File ID: GH099471A15Level: (low/med) LOW Date Received: 06/24/92% Moisture: 8 decanted: (Y/N) N Date Extracted: 06/26/92Concentrated Extract Volume: 500.0 (uL) Date Analyzed: 07/02/92Injection Volume: 2.0(uL) Dilution Factor: 1.0GPC Cleanup: (Y/N) Y pH: 6.1

CONCENTRATION UNITS:

(ug/L or ug/Kg) UG/KG Q

<u>108-95-2-----Phenol</u>	<u>360</u>	<u>U</u>
<u>111-44-4-----bis(2-Chloroethyl)Ether</u>	<u>360</u>	<u>U</u>
<u>95-57-8-----2-Chlorophenol</u>	<u>360</u>	<u>U</u>
<u>541-73-1-----1,3-Dichlorobenzene</u>	<u>360</u>	<u>U</u>
<u>106-46-7-----1,4-Dichlorobenzene</u>	<u>360</u>	<u>U</u>
<u>95-50-1-----1,2-Dichlorobenzene</u>	<u>360</u>	<u>U</u>
<u>95-48-7-----2-Methylphenol</u>	<u>360</u>	<u>U</u>
<u>108-60-1-----2,2'-Oxybis(1-Chloropropane)</u>	<u>360</u>	<u>U</u>
<u>106-44-5-----4-Methylphenol</u>	<u>360</u>	<u>U</u>
<u>621-64-7-----N-Nitroso-Di-n-Propylamine</u>	<u>360</u>	<u>U</u>
<u>67-72-1-----Hexachloroethane</u>	<u>360</u>	<u>U</u>
<u>98-95-3-----Nitrobenzene</u>	<u>360</u>	<u>U</u>
<u>78-59-1-----Isophorone</u>	<u>360</u>	<u>U</u>
<u>88-75-5-----2-Nitrophenol</u>	<u>360</u>	<u>U</u>
<u>105-67-9-----2,4-Dimethylphenol</u>	<u>360</u>	<u>U</u>
<u>111-91-1-----bis(2-Chloroethoxy)Methane</u>	<u>360</u>	<u>U</u>
<u>120-83-2-----2,4-Dichlorophenol</u>	<u>360</u>	<u>U</u>
<u>120-82-1-----1,2,4-Trichlorobenzene</u>	<u>360</u>	<u>U</u>
<u>91-20-3-----Naphthalene</u>	<u>360</u>	<u>U</u>
<u>106-47-8-----4-Chloroaniline</u>	<u>360</u>	<u>U</u>
<u>87-68-3-----Hexachlorobutadiene</u>	<u>360</u>	<u>U</u>
<u>59-50-7-----4-Chloro-3-Methylphenol</u>	<u>360</u>	<u>U</u>
<u>91-57-6-----2-Methylnaphthalene</u>	<u>360</u>	<u>U</u>
<u>77-47-4-----Hexachlorocyclopentadiene</u>	<u>360</u>	<u>U</u>
<u>88-06-2-----2,4,6-Trichlorophenol</u>	<u>360</u>	<u>U</u>
<u>95-95-4-----2,4,5-Trichlorophenol</u>	<u>870</u>	<u>U</u>
<u>91-58-7-----2-Choronaphthalene</u>	<u>360</u>	<u>U</u>
<u>88-74-4-----2-Nitroaniline</u>	<u>870</u>	<u>U</u>
<u>131-11-3-----Dimethyl Phthalate</u>	<u>360</u>	<u>U</u>
<u>208-96-8-----Acenaphthylene</u>	<u>360</u>	<u>U</u>
<u>606-20-2-----2,6-Dinitrotoluene</u>	<u>360</u>	<u>U</u>
<u>99-09-2-----3-Nitroaniline</u>	<u>870</u>	<u>U</u>
<u>83-32-9-----Acenaphthene</u>	<u>360</u>	<u>U</u>

1C
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEETORIGINAL
EPA SAMPLE NO.

CKY37

Lab Name: COMPUCHEM, RTPContract: 68D10083Lab Code: COMPUCase No.: 18347

SAS No.: _____

SDG No.: CKY28Matrix: (soil/water) SOILLab Sample ID: 499471Sample wt/vol: 30.0 (g/mL) GLab File ID: GH099471A15Level: (low/med) LOWDate Received: 06/24/92% Moisture: 8 decanted: (Y/N) NDate Extracted: 06/26/92Concentrated Extract Volume: 500.0 (uL)Date Analyzed: 07/02/92Injection Volume: 2.0 (uL)Dilution Factor: 1.0GPC Cleanup: (Y/N) Y pH: 6.1CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG

Q

CAS NO.	COMPOUND	Q
51-28-5-----	2,4-Dinitrophenol	870 U
100-02-7-----	4-Nitrophenol	870 U
132-64-9-----	Dibenzofuran	360 U
121-14-2-----	2,4-Dinitrotoluene	360 U
84-66-2-----	Diethylphthalate	360 U
7005-72-3-----	4-Chlorophenyl-phenylether	360 U
86-73-7-----	Fluorene	360 U
100-01-6-----	4-Nitroaniline	870 U
534-52-1-----	4,6-Dinitro-2-Methylphenol	870 U
86-30-6-----	N-Nitrosodiphenylamine (1)	360 U
101-55-3-----	4-Bromophenyl-phenylether	360 U
118-74-1-----	Hexachlorobenzene	360 U
87-86-5-----	Pentachlorophenol	870 U
85-01-8-----	Phenanthrene	360 U
120-12-7-----	Anthracene	360 U
86-74-8-----	Carbazole	360 U
84-74-2-----	Di-n-Butylphthalate	360 U
206-44-0-----	Fluoranthene	360 U
129-00-0-----	Pyrene	360 U
85-68-7-----	Butylbenzylphthalate	130 BJ
91-94-1-----	3,3'-Dichlorobenzidine	360 U
56-55-3-----	Benzo(a)Anthracene	360 U
218-01-9-----	Chrysene	360 U
117-81-7-----	bis(2-Ethylhexyl) Phthalate	360 U
117-84-0-----	Di-n-Octyl Phthalate	360 U
205-99-2-----	Benzo(b) Fluoranthene	360 U
207-08-9-----	Benzo(k) Fluoranthene	360 U
50-32-8-----	Benzo(a) Pyrene	360 U
193-39-5-----	Indeno(1,2,3-cd) Pyrene	360 U
53-70-3-----	Dibenz(a,h) Anthracene	360 U
191-24-2-----	Benzo(g,h,i) Perylene	360 U

(1) - Cannot be separated from Diphenylamine

1B
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

ORIGINAL
EPA SAMPLE NO.

CKY38

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 18347

SAS No.: _____

SDG No.: CKY28

Matrix: (soil/water) SOIL

Lab Sample ID: 499473

Sample wt/vol: 30.0 (g/mL) G

Lab File ID: GR099473A57

Level: (low/med) LOW

Date Received: 06/24/92

Moisture: 14 decanted: (Y/N) N

Date Extracted: 07/07/92

Concentrated Extract Volume: 500.0 (uL)

Date Analyzed: 07/09/92

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 6.4

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG

Q

CAS NO.	COMPOUND		
108-95-2-----	Phenol	380	U
111-44-4-----	bis(2-Chloroethyl)Ether	380	U
95-57-8-----	2-Chlorophenol	380	U
541-73-1-----	1,3-Dichlorobenzene	380	U
106-46-7-----	1,4-Dichlorobenzene	380	U
95-50-1-----	1,2-Dichlorobenzene	380	U
95-48-7-----	2-Methylphenol	380	U
108-60-1-----	2,2'-Oxybis(1-Chloropropane)	380	U
106-44-5-----	4-Methylphenol	380	U
621-64-7-----	N-Nitroso-Di-n-Propylamine	380	U
67-72-1-----	Hexachloroethane	380	U
98-95-3-----	Nitrobenzene	380	U
78-59-1-----	Isophorone	380	U
88-75-5-----	2-Nitrophenol	380	U
105-67-9-----	2,4-Dimethylphenol	380	U
111-91-1-----	bis(2-Chloroethoxy)Methane	380	U
120-83-2-----	2,4-Dichlorophenol	380	U
120-82-1-----	1,2,4-Trichlorobenzene	380	U
91-20-3-----	Naphthalene	380	U
106-47-8-----	4-Chloroaniline	380	U
87-68-3-----	Hexachlorobutadiene	380	U
59-50-7-----	4-Chloro-3-Methylphenol	380	U
91-57-6-----	2-Methylnaphthalene	380	U
77-47-4-----	Hexachlorocyclopentadiene	380	U
88-06-2-----	2,4,6-Trichlorophenol	380	U
95-95-4-----	2,4,5-Trichlorophenol	930	U
91-58-7-----	2-Chloronaphthalene	380	U
88-74-4-----	2-Nitroaniline	930	U
131-11-3-----	Dimethyl Phthalate	380	U
208-96-8-----	Acenaphthylene	380	U
606-20-2-----	2,6-Dinitrotoluene	380	U
99-09-2-----	3-Nitroaniline	930	U
83-32-9-----	Acenaphthene	380	U

1C
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

OPERA SAMPLE NO.

ORIGINAL
(Rev)

CKY38

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 18347

SAS No.: _____ SDG No.: CKY28

Matrix: (soil/water) SOIL

Lab Sample ID: 499473

Sample wt/vol: 30.0 (g/mL) G

Lab File ID: GR099473A57

Level: (low/med) LOW

Date Received: 06/24/92

% Moisture: 14 decanted: (Y/N) N

Date Extracted: 07/07/92

Concentrated Extract Volume: 500.0 (uL)

Date Analyzed: 07/09/92

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 6.4

CONCENTRATION UNITS:

(ug/L or ug/Kg) UG/KG

Q

CAS NO.	COMPOUND			
51-28-5-----	2,4-Dinitrophenol	930	U	
100-02-7-----	4-Nitrophenol	930	U	
132-64-9-----	Dibenzofuran	380	U	
121-14-2-----	2,4-Dinitrotoluene	380	U	
84-66-2-----	Diethylphthalate	380	U	
7005-72-3-----	4-Chlorophenyl-phenylether	380	U	
86-73-7-----	Fluorene	380	U	
100-01-6-----	4-Nitroaniline	930	U	
534-52-1-----	4,6-Dinitro-2-Methylphenol	930	U	
86-30-6-----	N-Nitrosodiphenylamine (1)	380	U	
101-55-3-----	4-Bromophenyl-phenylether	380	U	
118-74-1-----	Hexachlorobenzene	380	U	
87-86-5-----	Pentachlorophenol	930	U	
85-01-8-----	Phenanthrene	380	U	
120-12-7-----	Anthracene	380	U	
86-74-8-----	Carbazole	380	U	
84-74-2-----	Di-n-Butylphthalate	380	U	
206-44-0-----	Fluoranthene	61	J	
129-00-0-----	Pyrene	55	J	
85-68-7-----	Butylbenzylphthalate	380	U	
91-94-1-----	3,3'-Dichlorobenzidine	380	U	
56-55-3-----	Benzo(a)Anthracene	380	U	
218-01-9-----	Chrysene	380	U	
117-81-7-----	bis(2-Ethylhexyl)Phthalate	380	U	
117-84-0-----	Di-n-Octyl Phthalate	380	U	
205-99-2-----	Benzo(b)Fluoranthene	61	JX	
207-08-9-----	Benzo(k)Fluoranthene	61	JX	
50-32-8-----	Benzo(a)Pyrene	380	U	
193-39-5-----	Indeno(1,2,3-cd)Pyrene	380	U	
53-70-3-----	Dibenz(a,h)Anthracene	380	U	
191-24-2-----	Benzo(g,h,i)Perylene	380	U	

(1) - Cannot be separated from Diphenylamine

1D
PESTICIDE ORGANICS ANALYSIS DATA SHEETEPA SAMPLE NO.
(Med)

CKY14

Lab Name: COMPUCHEM RTPContract: 68D10083Lab Code: COMPU Case No.: 18347 SAS No.:SDG No.: CKY14Matrix: (soil/water) WATERLab Sample ID: 499425Sample wt/vol: 1000(g/ml)ML

Lab File ID:

% Moisture: decanted: (Y/N)

Date Received: 06/24/92Extraction: (SepF/Cont/Sonc) SEPFDate Extracted: 06/26/92Concentrated Extract Volume: 10000(uL)Date Analyzed: 06/30/92Injection Volume: 2.0(uL)Dilution Factor: 1GPC Cleanup: (Y/N) N

pH:

Sulfur Cleanup: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/L</u>	Q
319-84-6-----	alpha-BHC	0.050	U
319-85-7-----	beta-BHC	0.050	U
319-86-8-----	delta-BHC	0.050	U
58-89-9-----	gamma-BHC (Lindane)	0.050	U
76-44-8-----	Heptachlor	0.050	U
309-00-2-----	Aldrin	0.050	U
1024-57-3-----	Heptachlor epoxide	0.050	U
959-98-8-----	Endosulfan I	0.050	U
60-57-1-----	Dieldrin	0.10	U
72-55-9-----	4,4'-DDE	0.10	U
72-20-8-----	Endrin	0.10	U
33213-65-9-----	Endosulfan II	0.10	U
72-54-8-----	4,4'-DDD	0.10	U
1031-07-8-----	Endosulfan sulfate	0.10	U
50-29-3-----	4,4'-DDT	0.10	U
72-43-5-----	Methoxychlor	0.50	U
53494-70-5-----	Endrin ketone	0.10	U
7421-93-4-----	Endrin aldehyde	0.10	U
5103-71-9-----	alpha-Chlordane	0.050	U
5103-74-2-----	gamma-Chlordane	0.050	U
8001-35-2-----	Toxaphene	5.0	U
12674-11-2-----	Aroclor-1016	1.0	U
11104-28-2-----	Aroclor-1221	2.0	U
11141-16-5-----	Aroclor-1232	1.0	U
53469-21-9-----	Aroclor-1242	1.0	U
12672-29-6-----	Aroclor-1248	1.0	U
11097-69-1-----	Aroclor-1254	1.0	U
11096-82-5-----	Aroclor-1260	1.0	U

1D
PESTICIDE ORGANICS ANALYSIS DATA SHEET

EPA ⁰² SAMPLE NO.
(Ready)

CKY15

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 18347 SAS No.:

SDG No.: CKY14

Matrix: (soil/water) WATER

Lab Sample ID: 499435

Sample wt/vol: 1000(g/ml)ML

Lab File ID:

% Moisture: decanted: (Y/N)

Date Received: 06/24/92

Extraction: (SepF/Cont/Sonc) SEPF

Date Extracted: 06/26/92

Concentrated Extract Volume: 10000(uL)

Date Analyzed: 06/30/92

Injection Volume: 2.0(uL)

Dilution Factor: 1

GPC Cleanup: (Y/N) N

pH:

Sulfur Cleanup: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) UG/L	Q
319-84-6-----	alpha-BHC	0.050	U
319-85-7-----	beta-BHC	0.050	U
319-86-8-----	delta-BHC	0.050	U
58-89-9-----	gamma-BHC (Lindane)	0.050	U
76-44-8-----	Heptachlor	0.050	U
309-00-2-----	Aldrin	0.050	U
1024-57-3-----	Heptachlor epoxide	0.050	U
959-98-8-----	Endosulfan I	0.050	U
60-57-1-----	Dieldrin	0.10	U
72-55-9-----	4,4'-DDE	0.10	U
72-20-8-----	Endrin	0.10	U
33213-65-9-----	Endosulfan II	0.10	U
72-54-8-----	4,4'-DDD	0.10	U
1031-07-8-----	Endosulfan sulfate	0.10	U
50-29-3-----	4,4'-DDT	0.10	U
72-43-5-----	Methoxychlor	0.50	U
53494-70-5-----	Endrin ketone	0.10	U
7421-93-4-----	Endrin aldehyde	0.10	U
5103-71-9-----	alpha-Chlordane	0.050	U
5103-74-2-----	gamma-Chlordane	0.050	U
8001-35-2-----	Toxaphene	5.0	U
12674-11-2-----	Aroclor-1016	1.0	U
11104-28-2-----	Aroclor-1221	2.0	U
11141-16-5-----	Aroclor-1232	1.0	U
53469-21-9-----	Aroclor-1242	1.0	U
12672-29-6-----	Aroclor-1248	1.0	U
11097-69-1-----	Aroclor-1254	1.0	U
11096-82-5-----	Aroclor-1260	1.0	U

1D
PESTICIDE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.
(Ref)

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

CKY16

Lab Code: COMPU Case No.: 18347 SAS No.:

SDG No.: CKY14

Matrix: (soil/water) WATER

Lab Sample ID: 499436

Sample wt/vol: 1000(g/ml)ML

Lab File ID:

% Moisture: decanted: (Y/N)

Date Received: 06/24/92

Extraction: (SepF/Cont/Sonc) SEPF

Date Extracted: 06/26/92

Concentrated Extract Volume: 10000(uL)

Date Analyzed: 06/30/92

Injection Volume: 2.0(uL)

Dilution Factor: 1

GPC Cleanup: (Y/N) N pH:

Sulfur Cleanup: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) UG/L	Q
319-84-6-----	alpha-BHC	0.050	U
319-85-7-----	beta-BHC	0.050	U
319-86-8-----	delta-BHC	0.050	U
58-89-9-----	gamma-BHC (Lindane)	0.050	U
76-44-8-----	Heptachlor	0.050	U
309-00-2-----	Aldrin	0.050	U
1024-57-3-----	Heptachlor epoxide	0.050	U
959-98-8-----	Endosulfan I	0.050	U
60-57-1-----	Dieldrin	0.10	U
72-55-9-----	4,4'-DDE	0.10	U
72-20-8-----	Endrin	0.10	U
33213-65-9-----	Endosulfan II	0.10	U
72-54-8-----	4,4'-DDD	0.10	U
1031-07-8-----	Endosulfan sulfate	0.10	U
50-29-3-----	4,4'-DDT	0.10	U
72-43-5-----	Methoxychlor	0.50	U
53494-70-5-----	Endrin ketone	0.10	U
7421-93-4-----	Endrin aldehyde	0.10	U
5103-71-9-----	alpha-Chlordane	0.050	U
5103-74-2-----	gamma-Chlordane	0.050	U
8001-35-2-----	Toxaphene	5.0	U
12674-11-2-----	Aroclor-1016	1.0	U
11104-28-2-----	Aroclor-1221	2.0	U
11141-16-5-----	Aroclor-1232	1.0	U
53469-21-9-----	Aroclor-1242	1.0	U
12672-29-6-----	Aroclor-1248	1.0	U
11097-69-1-----	Aroclor-1254	1.0	U
11096-82-5-----	Aroclor-1260	1.0	U

1D
PESTICIDE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.
(Ready)

CKY17

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 18347 SAS No.:

SDG No.: CKY14

Matrix: (soil/water) WATER

Lab Sample ID: 499437

Sample wt/vol: 1000(g/ml) ML

Lab File ID:

% Moisture: decanted: (Y/N)

Date Received: 06/24/92

Extraction: (SepF/Cont/Sonc) SEPF

Date Extracted: 06/26/92

Concentrated Extract Volume: 10000(uL)

Date Analyzed: 06/30/92

Injection Volume: 2.0(uL)

Dilution Factor: 1

GPC Cleanup: (Y/N) N pH:

Sulfur Cleanup: (Y/N) N

CONCENTRATION UNITS:

(ug/L or ug/Kg) UG/L

Q

319-84-6-----	alpha-BHC	0.050	U
319-85-7-----	beta-BHC	0.050	U
319-86-8-----	delta-BHC	0.050	U
58-89-9-----	gamma-BHC (Lindane)	0.050	U
76-44-8-----	Heptachlor	0.0039	JP
309-00-2-----	Aldrin	0.050	U
1024-57-3-----	Heptachlor epoxide	0.050	U
959-98-8-----	Endosulfan I	0.050	U
60-57-1-----	Dieldrin	0.10	U
72-55-9-----	4,4'-DDE	0.10	U
72-20-8-----	Endrin	0.10	U
33213-65-9-----	Endosulfan II	0.10	U
72-54-8-----	4,4'-DDD	0.10	U
1031-07-8-----	Endosulfan sulfate	0.10	U
50-29-3-----	4,4'-DDT	0.10	U
72-43-5-----	Methoxychlor	0.50	U
53494-70-5-----	Endrin ketone	0.10	U
7421-93-4-----	Endrin aldehyde	0.10	U
5103-71-9-----	alpha-Chlordane	0.050	U
5103-74-2-----	gamma-Chlordane	0.050	U
8001-35-2-----	Toxaphene	5.0	U
12674-11-2-----	Aroclor-1016	1.0	U
11104-28-2-----	Aroclor-1221	2.0	U
11141-16-5-----	Aroclor-1232	1.0	U
53469-21-9-----	Aroclor-1242	1.0	U
12672-29-6-----	Aroclor-1248	1.0	U
11097-69-1-----	Aroclor-1254	1.0	U
11096-82-5-----	Aroclor-1260	1.0	U

1D
PESTICIDE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.
Original

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

CKY18

Lab Code: COMPU Case No.: 18347 SAS No.:

SDG No.: CKY14

Matrix: (soil/water) WATER

Lab Sample ID: 499439

Sample wt/vol: 1000(g/ml)ML

Lab File ID:

% Moisture: decanted: (Y/N)

Date Received: 06/24/92

Extraction: (SepF/Cont/Sonc) SEPF

Date Extracted: 06/26/92

Concentrated Extract Volume: 10000(uL)

Date Analyzed: 06/30/92

Injection Volume: 2.0(uL)

Dilution Factor: 1

GPC Cleanup: (Y/N) N

pH:

Sulfur Cleanup: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/L</u>	Q
319-84-6-----	alpha-BHC	0.050	U
319-85-7-----	beta-BHC	0.050	U
319-86-8-----	delta-BHC	0.050	U
58-89-9-----	gamma-BHC (Lindane)	0.0039	JP
76-44-8-----	Heptachlor	0.050	U
309-00-2-----	Aldrin	0.050	U
1024-57-3-----	Heptachlor epoxide	0.050	U
959-98-8-----	Endosulfan I	0.050	U
60-57-1-----	Dieldrin	0.10	U
72-55-9-----	4,4'-DDE	0.10	U
72-20-8-----	Endrin	0.10	U
33213-65-9-----	Endosulfan II	0.10	U
72-54-8-----	4,4'-DDD	0.10	U
1031-07-8-----	Endosulfan sulfate	0.10	U
50-29-3-----	4,4'-DDT	0.10	U
72-43-5-----	Methoxychlor	0.50	U
53494-70-5-----	Endrin ketone	0.10	U
7421-93-4-----	Endrin aldehyde	0.10	U
5103-71-9-----	alpha-Chlordane	0.050	U
5103-74-2-----	gamma-Chlordane	0.050	U
8001-35-2-----	Toxaphene	5.0	U
12674-11-2-----	Aroclor-1016	1.0	U
11104-28-2-----	Aroclor-1221	2.0	U
11141-16-5-----	Aroclor-1232	1.0	U
53469-21-9-----	Aroclor-1242	1.0	U
12672-29-6-----	Aroclor-1248	1.0	U
11097-69-1-----	Aroclor-1254	1.0	U
11096-82-5-----	Aroclor-1260	1.0	U

1D
PESTICIDE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

ORIGINAL
CKY19Lab Name: COMPUCHEM.RTPContract: 68D10083

Lab Code: COMPU Case No.: 18347 SAS No.: SDG No.: CKY14
 Matrix: (soil/water) WATER Lab Sample ID: 499440
 Sample wt/vol: 1000(g/ml)ML Lab File ID:
 % Moisture: decanted: (Y/N) Date Received: 06/24/92
 Extraction: (SepF/Cont/Sonc) SEPF Date Extracted: 06/26/92
 Concentrated Extract Volume: 10000(uL) Date Analyzed: 06/30/92
 Injection Volume: 2.0(uL) Dilution Factor: 1
 GPC Cleanup: (Y/N) N Sulfur Cleanup: (Y/N) N

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/LQ

319-84-6-----	alpha-BHC	0.050	U
319-85-7-----	beta-BHC	0.050	U
319-86-8-----	delta-BHC	0.050	U
58-89-9-----	gamma-BHC (Lindane)	0.050	U
76-44-8-----	Heptachlor	0.050	U
309-00-2-----	Aldrin	0.050	U
1024-57-3-----	Heptachlor epoxide	0.050	U
959-98-8-----	Endosulfan I	0.050	U
60-57-1-----	Dieldrin	0.10	U
72-55-9-----	4,4'-DDE	0.10	U
72-20-8-----	Endrin	0.10	U
33213-65-9-----	Endosulfan II	0.10	U
72-54-8-----	4,4'-DDD	0.10	U
1031-07-8-----	Endosulfan sulfate	0.10	U
50-29-3-----	4,4'-DDT	0.10	U
72-43-5-----	Methoxychlor	0.50	U
53494-70-5-----	Endrin ketone	0.10	U
7421-93-4-----	Endrin aldehyde	0.10	U
5103-71-9-----	alpha-Chlordane	0.050	U
5103-74-2-----	gamma-Chlordane	0.050	U
8001-35-2-----	Toxaphene	5.0	U
12674-11-2-----	Aroclor-1016	1.0	U
11104-28-2-----	Aroclor-1221	2.0	U
11141-16-5-----	Aroclor-1232	1.0	U
53469-21-9-----	Aroclor-1242	1.0	U
12672-29-6-----	Aroclor-1248	1.0	U
11097-69-1-----	Aroclor-1254	1.0	U
11096-82-5-----	Aroclor-1260	1.0	U

1D
PESTICIDE ORGANICS ANALYSIS DATA SHEETEPA SAMPLE NO.
*021
Med*Lab Name: COMPUCHEM, RTPContract: 68D10083CKY20Lab Code: COMPU Case No.: 18347 SAS No.: SDG No.: CKY14Matrix: (soil/water) WATERLab Sample ID: 499441Sample wt/vol: 1000(g/ml)ML

Lab File ID:

% Moisture: decanted: (Y/N)

Date Received: 06/24/92Extraction: (SepF/Cont/Sonc) SEPFDate Extracted: 06/26/92Concentrated Extract Volume: 10000(uL)Date Analyzed: 07/01/92Injection Volume: 2.0(uL)Dilution Factor: 1GPC Cleanup: (Y/N) N pH:Sulfur Cleanup: (Y/N) N

CONCENTRATION UNITS:

(ug/L or ug/Kg) UG/L

Q

CAS NO.	COMPOUND	Q
319-84-6-----	alpha-BHC	0.050 U
319-85-7-----	beta-BHC	0.050 U
319-86-8-----	delta-BHC	0.050 U
58-89-9-----	gamma-BHC (Lindane)	0.050 U
76-44-8-----	Heptachlor	0.0099 J
309-00-2-----	Aldrin	0.050 U
1024-57-3-----	Heptachlor epoxide	0.050 U
959-98-8-----	Endosulfan I	0.050 U
60-57-1-----	Dieldrin	0.10 U
72-55-9-----	4,4'-DDE	0.10 U
72-20-8-----	Endrin	0.10 U
33213-65-9-----	Endosulfan II	0.10 U
72-54-8-----	4,4'-DDD	0.10 U
1031-07-8-----	Endosulfan sulfate	0.10 U
50-29-3-----	4,4'-DDT	0.10 U
72-43-5-----	Methoxychlor	0.50 U
53494-70-5-----	Endrin ketone	0.10 U
7421-93-4-----	Endrin aldehyde	0.10 U
5103-71-9-----	alpha-Chlordane	0.050 U
5103-74-2-----	gamma-Chlordane	0.050 U
8001-35-2-----	Toxaphene	5.0 U
12674-11-2-----	Aroclor-1016	1.0 U
11104-28-2-----	Aroclor-1221	2.0 U
11141-16-5-----	Aroclor-1232	1.0 U
53469-21-9-----	Aroclor-1242	1.0 U
12672-29-6-----	Aroclor-1248	1.0 U
11097-69-1-----	Aroclor-1254	1.0 U
11096-82-5-----	Aroclor-1260	1.0 U

1D
PESTICIDE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

(RSD)

CKY22

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 18347 SAS No.:

SDG No.: CKY14

Matrix: (soil/water) WATER

Lab Sample ID: 499442

Sample wt/vol: 1000(g/ml)ML

Lab File ID:

Moisture: decanted: (Y/N)

Date Received: 06/24/92

Extraction: (SepF/Cont/Sonc) SEPF

Date Extracted: 06/26/92

Concentrated Extract Volume: 10000(uL)

Date Analyzed: 07/01/92

Injection Volume: 2.0(uL)

Dilution Factor: 1

GPC Cleanup: (Y/N) N pH:

Sulfur Cleanup: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/L</u>	Q
319-84-6-----	alpha-BHC	0.050	U
319-85-7-----	beta-BHC	0.050	U
319-86-8-----	delta-BHC	0.050	U
58-89-9-----	gamma-BHC (Lindane)	0.050	U
76-44-8-----	Heptachlor	0.050	U
309-00-2-----	Aldrin	0.050	U
1024-57-3-----	Heptachlor epoxide	0.050	U
959-98-8-----	Endosulfan I	0.050	U
60-57-1-----	Die�drin	0.10	U
72-55-9-----	4,4'-DDE	0.10	U
72-20-8-----	Endrin	0.10	U
33213-65-9-----	Endosulfan II	0.10	U
72-54-8-----	4,4'-DDD	0.10	U
1031-07-8-----	Endosulfan sulfate	0.10	U
50-29-3-----	4,4'-DDT	0.10	U
72-43-5-----	Methoxychlor	0.50	U
53494-70-5-----	Endrin ketone	0.10	U
7421-93-4-----	Endrin aldehyde	0.10	U
5103-71-9-----	alpha-Chlordane	0.050	U
5103-74-2-----	gamma-Chlordane	0.050	U
8001-35-2-----	Toxaphene	5.0	U
12674-11-2-----	Aroclor-1016	1.0	U
11104-28-2-----	Aroclor-1221	2.0	U
11141-16-5-----	Aroclor-1232	1.0	U
53469-21-9-----	Aroclor-1242	1.0	U
12672-29-6-----	Aroclor-1248	1.0	U
11097-69-1-----	Aroclor-1254	1.0	U
11096-82-5-----	Aroclor-1260	1.0	U

1D
PESTICIDE ORGANICS ANALYSIS DATA SHEET

EPAOR SAMPLE NO.

ORIGINAL

Mod

CKY23

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 18347 SAS No.:

SDG No.: CKY14

Matrix: (soil/water) WATER

Lab Sample ID: 499444

Sample wt/vol: 1000(g/ml)ML

Lab File ID:

% Moisture: decanted: (Y/N)

Date Received: 06/24/92

Extraction: (SepF/Cont/Sonc) SEPF

Date Extracted: 06/26/92

Concentrated Extract Volume: 10000(uL)

Date Analyzed: 07/01/92

Injection Volume: 2.0(uL)

Dilution Factor: 1

GPC Cleanup: (Y/N) N pH:

Sulfur Cleanup: (Y/N) N

CONCENTRATION UNITS:

(ug/L or ug/Kg) UG/L

Q

319-84-6-----	alpha-BHC	0.050	U
319-85-7-----	beta-BHC	0.050	U
319-86-8-----	delta-BHC	0.050	U
58-89-9-----	gamma-BHC (Lindane)	0.050	U
76-44-8-----	Heptachlor	0.050	U
309-00-2-----	Aldrin	0.050	U
1024-57-3-----	Heptachlor epoxide	0.050	U
959-98-8-----	Endosulfan I	0.050	U
60-57-1-----	Dieldrin	0.10	U
72-55-9-----	4,4'-DDE	0.10	U
72-20-8-----	Endrin	0.10	U
33213-65-9-----	Endosulfan II	0.10	U
72-54-8-----	4,4'-DDD	0.10	U
1031-07-8-----	Endosulfan sulfate	0.10	U
50-29-3-----	4,4'-DDT	0.10	U
72-43-5-----	Methoxychlor	0.50	U
53494-70-5-----	Endrin ketone	0.10	U
7421-93-4-----	Endrin aldehyde	0.10	U
5103-71-9-----	alpha-Chlordane	0.050	U
5103-74-2-----	gamma-Chlordane	0.050	U
8001-35-2-----	Toxaphene	5.0	U
12674-11-2-----	Aroclor-1016	1.0	U
11104-28-2-----	Aroclor-1221	2.0	U
11141-16-5-----	Aroclor-1232	1.0	U
53469-21-9-----	Aroclor-1242	1.0	U
12672-29-6-----	Aroclor-1248	1.0	U
11097-69-1-----	Aroclor-1254	1.0	U
11096-82-5-----	Aroclor-1260	1.0	U

1D
PESTICIDE ORGANICS ANALYSIS DATA SHEET

ORIGIN
EPA SAMPLE NO.
(Ind)

CKY24

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 18347 SAS No.:

SDG No.: CKY14

Matrix: (soil/water) WATER

Lab Sample ID: 499445

Sample wt/vol: 1000(g/ml)ML

Lab File ID:

% Moisture: decanted: (Y/N)

Date Received: 06/24/92

Extraction: (SepF/Cont/Sonc) SEPF

Date Extracted: 06/26/92

Concentrated Extract Volume: 10000(uL)

Date Analyzed: 07/01/92

Injection Volume: 2.0(uL)

Dilution Factor: 1

GPC Cleanup: (Y/N) N

pH:

Sulfur Cleanup: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/L</u>	Q
319-84-6-----	alpha-BHC	0.050	U
319-85-7-----	beta-BHC	0.050	U
319-86-8-----	delta-BHC	0.050	U
58-89-9-----	gamma-BHC (Lindane)	0.050	U
76-44-8-----	Heptachlor	0.050	U
309-00-2-----	Aldrin	0.050	U
1024-57-3-----	Heptachlor epoxide	0.050	U
959-98-8-----	Endosulfan I	0.050	U
60-57-1-----	Dieldrin	0.10	U
72-55-9-----	4,4'-DDE	0.10	U
72-20-8-----	Endrin	0.10	U
33213-65-9-----	Endosulfan II	0.10	U
72-54-8-----	4,4'-DDD	0.10	U
1031-07-8-----	Endosulfan sulfate	0.10	U
50-29-3-----	4,4'-DDT	0.10	U
72-43-5-----	Methoxychlor	0.50	U
53494-70-5-----	Endrin ketone	0.10	U
7421-93-4-----	Endrin aldehyde	0.10	U
5103-71-9-----	alpha-Chlordane	0.050	U
5103-74-2-----	gamma-Chlordane	0.050	U
8001-35-2-----	Toxaphene	5.0	U
12674-11-2-----	Aroclor-1016	1.0	U
11104-28-2-----	Aroclor-1221	2.0	U
11141-16-5-----	Aroclor-1232	1.0	U
53469-21-9-----	Aroclor-1242	1.0	U
12672-29-6-----	Aroclor-1248	1.0	U
11097-69-1-----	Aroclor-1254	1.0	U
11096-82-5-----	Aroclor-1260	1.0	U

1D
PESTICIDE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

(Ready)

CKY25

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 18347 SAS No.:

SDG No.: CKY14

Matrix: (soil/water) WATER

Lab Sample ID: 499447

Sample wt/vol: 1000(g/ml)ML

Lab File ID:

% Moisture: decanted: (Y/N)

Date Received: 06/24/92

Extraction: (SepF/Cont/Sonc) SEPF

Date Extracted: 06/26/92

Concentrated Extract Volume: 10000(uL)

Date Analyzed: 07/01/92

Injection Volume: 2.0(uL)

Dilution Factor: 1

GPC Cleanup: (Y/N) N pH:

Sulfur Cleanup: (Y/N) N

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/L

Q

CAS NO.	COMPOUND	CONCENTRATION UNITS:	
		(ug/L or ug/Kg) <u>UG/L</u>	<u>Q</u>
319-84-6-----	alpha-BHC	0.050	U
319-85-7-----	beta-BHC	0.050	U
319-86-8-----	delta-BHC	0.050	U
58-89-9-----	gamma-BHC (Lindane)	0.050	U
76-44-8-----	Heptachlor	0.050	U
309-00-2-----	Aldrin	0.050	U
1024-57-3-----	Heptachlor epoxide	0.050	U
959-98-8-----	Endosulfan I	0.050	U
60-57-1-----	Dieldrin	0.10	U
72-55-9-----	4,4'-DDE	0.10	U
72-20-8-----	Endrin	0.10	U
33213-65-9-----	Endosulfan II	0.10	U
72-54-8-----	4,4'-DDD	0.10	U
1031-07-8-----	Endosulfan sulfate	0.10	U
50-29-3-----	4,4'-DDT	0.10	U
72-43-5-----	Methoxychlor	0.50	U
53494-70-5-----	Endrin ketone	0.10	U
7421-93-4-----	Endrin aldehyde	0.10	U
5103-71-9-----	alpha-Chlordane	0.050	U
5103-74-2-----	gamma-Chlordane	0.050	U
8001-35-2-----	Toxaphene	5.0	U
12674-11-2-----	Aroclor-1016	1.0	U
11104-28-2-----	Aroclor-1221	2.0	U
11141-16-5-----	Aroclor-1232	1.0	U
53469-21-9-----	Aroclor-1242	1.0	U
12672-29-6-----	Aroclor-1248	1.0	U
11097-69-1-----	Aroclor-1254	1.0	U
11096-82-5-----	Aroclor-1260	1.0	U

1D
PESTICIDE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.
(Ready)

CKY26

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 18347 SAS No.:

SDG No.: CKY14

Matrix: (soil/water) WATER

Lab Sample ID: 499449

Sample wt/vol: 1000(g/ml)ML

Lab File ID:

% Moisture: decanted: (Y/N)

Date Received: 06/24/92

Extraction: (SepF/Cont/Sonc) SEPF

Date Extracted: 06/26/92

Concentrated Extract Volume: 10000(uL)

Date Analyzed: 07/01/92

Injection Volume: 2.0(uL)

Dilution Factor: 1

GPC Cleanup: (Y/N) N pH:

Sulfur Cleanup: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/L</u>	Q
---------	----------	---	---

319-84-6-----	alpha-BHC	0.050	U
319-85-7-----	beta-BHC	0.050	U
319-86-8-----	delta-BHC	0.050	U
58-89-9-----	gamma-BHC (Lindane)	0.050	U
76-44-8-----	Heptachlor	0.050	U
309-00-2-----	Aldrin	0.050	U
1024-57-3-----	Heptachlor epoxide	0.050	U
959-98-8-----	Endosulfan I	0.050	U
60-57-1-----	Dieldrin	0.10	U
72-55-9-----	4,4'-DDE	0.10	U
72-20-8-----	Endrin	0.10	U
33213-65-9-----	Endosulfan II	0.10	U
72-54-8-----	4,4'-DDD	0.10	U
1031-07-8-----	Endosulfan sulfate	0.0082	J
50-29-3-----	4,4'-DDT	0.10	U
72-43-5-----	Methoxychlor	0.019	JP
53494-70-5-----	Endrin ketone	0.10	U
7421-93-4-----	Endrin aldehyde	0.10	U
5103-71-9-----	alpha-Chlordane	0.050	U
5103-74-2-----	gamma-Chlordane	0.050	U
8001-35-2-----	Toxaphene	5.0	U
12674-11-2-----	Aroclor-1016	1.0	U
11104-28-2-----	Aroclor-1221	2.0	U
11141-16-5-----	Aroclor-1232	1.0	U
53469-21-9-----	Aroclor-1242	1.0	U
12672-29-6-----	Aroclor-1248	1.0	U
11097-69-1-----	Aroclor-1254	1.0	U
11096-82-5-----	Aroclor-1260	1.0	U

1D
PESTICIDE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

(Ready)

CKY28

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 18347 SAS No.:

SDG No.: CKY28

Matrix: (soil/water) SOIL

Lab Sample ID: 499452

Sample wt/vol: 30.00(g/ml)G

Lab File ID:

% Moisture: 19 decanted: (Y/N) N

Date Received: 06/24/92

Extraction: (SepF/Cont/Sonc) SONC

Date Extracted: 06/26/92

Concentrated Extract Volume: 5000(uL)

Date Analyzed: 07/06/92

Injection Volume: 2.0(uL)

Dilution Factor: 1

GPC Cleanup: (Y/N) Y pH: 7.2

Sulfur Cleanup: (Y/N) N

CONCENTRATION UNITS:

(ug/L or ug/Kg) UG/KG

Q

CAS NO.	COMPOUND	CONCENTRATION UNITS:	
		(ug/L or ug/Kg) <u>UG/KG</u>	Q
319-84-6-----	alpha-BHC	2.1	U
319-85-7-----	beta-BHC	2.1	U
319-86-8-----	delta-BHC	2.1	U
58-89-9-----	gamma-BHC (Lindane)	2.1	U
76-44-8-----	Heptachlor	0.17	JP
309-00-2-----	Aldrin	2.1	U
1024-57-3-----	Heptachlor epoxide	2.1	U
959-98-8-----	Endosulfan I	2.1	U
60-57-1-----	Dieldrin	4.1	U
72-55-9-----	4,4'-DDE	0.20	JP
72-20-8-----	Endrin	4.1	U
33213-65-9-----	Endosulfan II	4.1	U
72-54-8-----	4,4'-DDD	4.1	U
1031-07-8-----	Endosulfan sulfate	4.1	U
50-29-3-----	4,4'-DDT	4.1	U
72-43-5-----	Methoxychlor	21	U
53494-70-5-----	Endrin ketone	0.18	JP
7421-93-4-----	Endrin aldehyde	4.1	U
5103-71-9-----	alpha-Chlordane	2.1	U
5103-74-2-----	gamma-Chlordane	2.1	U
8001-35-2-----	Toxaphene	210	U
12674-11-2-----	Aroclor-1016	41	U
11104-28-2-----	Aroclor-1221	83	U
11141-16-5-----	Aroclor-1232	41	U
53469-21-9-----	Aroclor-1242	41	U
12672-29-6-----	Aroclor-1248	41	U
11097-69-1-----	Aroclor-1254	41	U
11096-82-5-----	Aroclor-1260	41	U

1D
PESTICIDE ORGANICS ANALYSIS DATA SHEET

ORIGINAL
EPA SAMPLE NO.

CKY29

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 18347 SAS No.:

SDG No.: CKY28

Matrix: (soil/water) SOIL

Lab Sample ID: 499459

Sample wt/vol: 30.00(g/ml)G

Lab File ID:

% Moisture: 17 decanted: (Y/N) N

Date Received: 06/24/92

Extraction: (SepF/Cont/Sonc) SONC

Date Extracted: 06/26/92

Concentrated Extract Volume: 5000(uL)

Date Analyzed: 07/07/92

Injection Volume: 2.0(uL)

Dilution Factor: 1

GPC Cleanup: (Y/N) Y pH: 6.6

Sulfur Cleanup: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/KG</u>	Q
319-84-6-----	alpha-BHC	2.0	U
319-85-7-----	beta-BHC	2.0	U
319-86-8-----	delta-BHC	2.0	U
58-89-9-----	gamma-BHC (Lindane)	0.079	JP
76-44-8-----	Heptachlor	0.35	JP
309-00-2-----	Aldrin	2.0	U
1024-57-3-----	Heptachlor epoxide	2.0	U
959-98-8-----	Endosulfan I	2.0	U
60-57-1-----	Dieldrin	0.072	JP
72-55-9-----	4,4'-DDE	0.64	JP
72-20-8-----	Endrin	0.47	JP
33213-65-9-----	Endosulfan II	4.0	U
72-54-8-----	4,4'-DDD	0.15	JP
1031-07-8-----	Endosulfan sulfate	4.0	U
50-29-3-----	4,4'-DDT	4.0	U
72-43-5-----	Methoxychlor	0.32	JPB
53494-70-5-----	Endrin ketone	0.31	JP
7421-93-4-----	Endrin aldehyde	4.0	U
5103-71-9-----	alpha-Chlordane	0.36	JP
5103-74-2-----	gamma-Chlordane	0.46	JP
8001-35-2-----	Toxaphene	200	U
12674-11-2-----	Aroclor-1016	40	U
11104-28-2-----	Aroclor-1221	81	U
11141-16-5-----	Aroclor-1232	40	U
53469-21-9-----	Aroclor-1242	40	U
12672-29-6-----	Aroclor-1248	40	U
11097-69-1-----	Aroclor-1254	40	U
11096-82-5-----	Aroclor-1260	40	U

1D
PESTICIDE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

(Ind)

CKY30

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 18347 SAS No.:

SDG No.: CKY28

Matrix: (soil/water) SOIL

Lab Sample ID: 499463

Sample wt/vol: 30.00(g/ml)G

Lab File ID:

% Moisture: 40 decanted: (Y/N) N

Date Received: 06/24/92

Extraction: (SepF/Cont/Sonc) SONC

Date Extracted: 06/26/92

Concentrated Extract Volume: 5000(uL)

Date Analyzed: 07/07/92

Injection Volume: 2.0(uL)

Dilution Factor: 1

GPC Cleanup: (Y/N) Y pH: 5.5

Sulfur Cleanup: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/KG</u>	Q
319-84-6-----	alpha-BHC	2.8	U
319-85-7-----	beta-BHC	2.8	U
319-86-8-----	delta-BHC	2.8	U
58-89-9-----	gamma-BHC (Lindane)	2.8	U
76-44-8-----	Heptachlor	0.48	JP
309-00-2-----	Aldrin	2.8	U
1024-57-3-----	Heptachlor epoxide	2.8	U
959-98-8-----	Endosulfan I	2.8	U
60-57-1-----	Dieldrin	5.5	U
72-55-9-----	4,4'-DDE	0.39	JP
72-20-8-----	Endrin	5.5	U
33213-65-9-----	Endosulfan II	5.5	U
72-54-8-----	4,4'-DDD	5.5	U
1031-07-8-----	Endosulfan sulfate	5.5	U
50-29-3-----	4,4'-DDT	5.5	U
72-43-5-----	Methoxychlor	28	U
53494-70-5-----	Endrin ketone	5.5	U
7421-93-4-----	Endrin aldehyde	5.5	U
5103-71-9-----	alpha-Chlordane	2.8	U
5103-74-2-----	gamma-Chlordane	2.8	U
8001-35-2-----	Toxaphene	280	U
12674-11-2-----	Aroclor-1016	55	U
11104-28-2-----	Aroclor-1221	110	U
11141-16-5-----	Aroclor-1232	55	U
53469-21-9-----	Aroclor-1242	55	U
12672-29-6-----	Aroclor-1248	55	U
11097-69-1-----	Aroclor-1254	55	U
11096-82-5-----	Aroclor-1260	55	U

1D
PESTICIDE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CKY31

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 18347 SAS No.:

SDG No.: CKY28

Matrix: (soil/water) SOIL

Lab Sample ID: 499464

Sample wt/vol: 30.00(g/ml)G

Lab File ID:

% Moisture: 23 decanted: (Y/N) N

Date Received: 06/24/92

Extraction: (SepF/Cont/Sonc) SONC

Date Extracted: 06/26/92

Concentrated Extract Volume: 5000(uL)

Date Analyzed: 07/07/92

Injection Volume: 2.0(uL)

Dilution Factor: 1

GPC Cleanup: (Y/N) Y pH: 7.1

Sulfur Cleanup: (Y/N) N

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG

CAS NO.	COMPOUND	Q
319-84-6-----	alpha-BHC	2.2 U
319-85-7-----	beta-BHC	2.2 U
319-86-8-----	delta-BHC	2.2 U
58-89-9-----	gamma-BHC (Lindane)	2.2 U
76-44-8-----	Heptachlor	0.25 JP
309-00-2-----	Aldrin	2.2 U
1024-57-3-----	Heptachlor epoxide	2.2 U
959-98-8-----	Endosulfan I	2.2 U
60-57-1-----	Dieldrin	4.3 U
72-55-9-----	4,4'-DDE	0.38 J
72-20-8-----	Endrin	4.3 U
33213-65-9-----	Endosulfan II	4.3 U
72-54-8-----	4,4'-DDD	0.51 JP
1031-07-8-----	Endosulfan sulfate	4.3 U
50-29-3-----	4,4'-DDT	0.15 JP
72-43-5-----	Methoxychlor	0.92 JPB
53494-70-5-----	Endrin ketone	4.3 U
7421-93-4-----	Endrin aldehyde	4.3 U
5103-71-9-----	alpha-Chlordane	2.2 U
5103-74-2-----	gamma-Chlordane	0.16 JP
8001-35-2-----	Toxaphene	220 U
12674-11-2-----	Aroclor-1016	43 U
11104-28-2-----	Aroclor-1221	87 U
11141-16-5-----	Aroclor-1232	43 U
53469-21-9-----	Aroclor-1242	43 U
12672-29-6-----	Aroclor-1248	43 U
11097-69-1-----	Aroclor-1254	43 U
11096-82-5-----	Aroclor-1260	43 U

1D
PESTICIDE ORGANICS ANALYSIS DATA SHEETORIG
EPA SAMPLE NO.
(red)

CKY32

Lab Name: COMPUCHEM, RTPContract: 68D10083Lab Code: COMPU Case No.: 18347 SAS No.:SDG No.: CKY28Matrix: (soil/water) SOILLab Sample ID: 499465Sample wt/vol: 30.00(g/ml)G

Lab File ID:

% Moisture: 18 decanted: (Y/N) NDate Received: 06/24/92Extraction: (SepF/Cont/Sonc) SONCDate Extracted: 06/26/92Concentrated Extract Volume: 5000(uL)Date Analyzed: 07/06/92Injection Volume: 2.0(uL)Dilution Factor: 1GPC Cleanup: (Y/N) Y pH: 6.5Sulfur Cleanup: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/KG</u>	Q
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319-84-6-----	alpha-BHC	2.1	U
319-85-7-----	beta-BHC	2.1	U
319-86-8-----	delta-BHC	2.1	U
58-89-9-----	gamma-BHC (Lindane)	0.25	JP
76-44-8-----	Heptachlor	0.13	JP
309-00-2-----	Aldrin	2.1	U
1024-57-3-----	Heptachlor epoxide	2.1	U
959-98-8-----	Endosulfan I	2.1	U
60-57-1-----	Dieldrin	0.20	JP
72-55-9-----	4,4'-DDE	4.0	U
72-20-8-----	Endrin	4.0	U
33213-65-9-----	Endosulfan II	4.0	U
72-54-8-----	4,4'-DDD	4.0	U
1031-07-8-----	Endosulfan sulfate	4.0	U
50-29-3-----	4,4'-DDT	0.16	JP
72-43-5-----	Methoxychlor	21	U
53494-70-5-----	Endrin ketone	0.66	JP
7421-93-4-----	Endrin aldehyde	4.0	U
5103-71-9-----	alpha-Chlordane	2.1	U
5103-74-2-----	gamma-Chlordane	0.072	JP
8001-35-2-----	Toxaphene	210	U
12674-11-2-----	Aroclor-1016	40	U
11104-28-2-----	Aroclor-1221	82	U
11141-16-5-----	Aroclor-1232	40	U
53469-21-9-----	Aroclor-1242	40	U
12672-29-6-----	Aroclor-1248	40	U
11097-69-1-----	Aroclor-1254	40	U
11096-82-5-----	Aroclor-1260	40	U

1D
PESTICIDE ORGANICS ANALYSIS DATA SHEET

ORIGINAL
(Red)
EPA SAMPLE NO.

CKY33

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 18347 SAS No.: SDG No.: CKY28

Matrix: (soil/water) SOIL

Lab Sample ID: 499466

Sample wt/vol: 30.00(g/ml)G

Lab File ID:

% Moisture: 11 decanted: (Y/N) N

Date Received: 06/24/92

Extraction: (SepF/Cont/Sonc) SONC

Date Extracted: 06/26/92

Concentrated Extract Volume: 5000(uL)

Date Analyzed: 07/07/92

Injection Volume: 2.0(uL)

Dilution Factor: 1

GPC Cleanup: (Y/N) Y pH: 5.9

Sulfur Cleanup: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/KG</u>	Q
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319-84-6-----	alpha-BHC	0.25	JP
319-85-7-----	beta-BHC	1.9	U
319-86-8-----	delta-BHC	1.9	U
58-89-9-----	gamma-BHC (Lindane)	0.16	JP
76-44-8-----	Heptachlor	0.065	JP
309-00-2-----	Aldrin	0.65	J
1024-57-3-----	Heptachlor epoxide	0.18	JP
959-98-8-----	Endosulfan I	0.13	JP
60-57-1-----	Dieldrin	3.7	U
72-55-9-----	4,4'-DDE	3.7	U
72-20-8-----	Endrin	0.075	JP
33213-65-9-----	Endosulfan II	3.7	U
72-54-8-----	4,4'-DDD	3.7	U
1031-07-8-----	Endosulfan sulfate	0.096	JP
50-29-3-----	4,4'-DDT	3.7	U
72-43-5-----	Methoxychlor	0.39	JB
53494-70-5-----	Endrin ketone	3.7	U
7421-93-4-----	Endrin aldehyde	3.7	U
5103-71-9-----	alpha-Chlordane	1.9	U
5103-74-2-----	gamma-Chlordane	0.41	JP
8001-35-2-----	Toxaphene	190	U
12674-11-2-----	Aroclor-1016	37	U
11104-28-2-----	Aroclor-1221	75	U
11141-16-5-----	Aroclor-1232	37	U
53469-21-9-----	Aroclor-1242	37	U
12672-29-6-----	Aroclor-1248	37	U
11097-69-1-----	Aroclor-1254	37	U
11096-82-5-----	Aroclor-1260	37	U

PESTICIDE ORGANICS ANALYSIS DATA SHEET

CKY34

Lab Name: COMPUCHEM, RTPContract: 68D10083Lab Code: COMPU Case No.: 18347 SAS No.: SDG No.: CKY28Matrix: (soil/water) SOILLab Sample ID: 499467Sample wt/vol: 30.00(g/ml)G

Lab File ID:

% Moisture: 13 decanted: (Y/N) NDate Received: 06/24/92Extraction: (SepF/Cont/Sonc) SONCDate Extracted: 06/26/92Concentrated Extract Volume: 5000(uL)Date Analyzed: 07/07/92Injection Volume: 2.0(uL)Dilution Factor: 1GPC Cleanup: (Y/N) Y pH: 6.5Sulfur Cleanup: (Y/N) NCONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG

Q

<u>319-84-6-----alpha-BHC</u>	<u>2.0</u>	<u>U</u>
<u>319-85-7-----beta-BHC</u>	<u>2.0</u>	<u>U</u>
<u>319-86-8-----delta-BHC</u>	<u>2.0</u>	<u>U</u>
<u>58-89-9-----gamma-BHC (Lindane)</u>	<u>0.13</u>	<u>JP</u>
<u>76-44-8-----Heptachlor</u>	<u>2.0</u>	<u>U</u>
<u>309-00-2-----Aldrin</u>	<u>0.58</u>	<u>JP</u>
<u>1024-57-3-----Heptachlor epoxide</u>	<u>0.30</u>	<u>JP</u>
<u>959-98-8-----Endosulfan I</u>	<u>0.40</u>	<u>J</u>
<u>60-57-1-----Dieldrin</u>	<u>3.8</u>	<u>U</u>
<u>72-55-9-----4,4'-DDE</u>	<u>3.8</u>	<u>U</u>
<u>72-20-8-----Endrin</u>	<u>0.59</u>	<u>JP</u>
<u>33213-65-9-----Endosulfan II</u>	<u>0.31</u>	<u>JP</u>
<u>72-54-8-----4,4'-DDD</u>	<u>3.8</u>	<u>U</u>
<u>1031-07-8-----Endosulfan sulfate</u>	<u>3.8</u>	<u>U</u>
<u>50-29-3-----4,4'-DDT</u>	<u>3.8</u>	<u>U</u>
<u>72-43-5-----Methoxychlor</u>	<u>0.46</u>	<u>JPB</u>
<u>53494-70-5-----Endrin ketone</u>	<u>0.15</u>	<u>JP</u>
<u>7421-93-4-----Endrin aldehyde</u>	<u>3.8</u>	<u>U</u>
<u>5103-71-9-----alpha-Chlordane</u>	<u>0.55</u>	<u>JP</u>
<u>5103-74-2-----gamma-Chlordane</u>	<u>0.67</u>	<u>JP</u>
<u>8001-35-2-----Toxaphene</u>	<u>200</u>	<u>U</u>
<u>12674-11-2-----Aroclor-1016</u>	<u>38</u>	<u>U</u>
<u>11104-28-2-----Aroclor-1221</u>	<u>77</u>	<u>U</u>
<u>11141-16-5-----Aroclor-1232</u>	<u>38</u>	<u>U</u>
<u>53469-21-9-----Aroclor-1242</u>	<u>38</u>	<u>U</u>
<u>12672-29-6-----Aroclor-1248</u>	<u>38</u>	<u>U</u>
<u>11097-69-1-----Aroclor-1254</u>	<u>38</u>	<u>U</u>
<u>11096-82-5-----Aroclor-1260</u>	<u>38</u>	<u>U</u>

ID
PESTICIDE ORGANICS ANALYSIS DATA SHEET

ORIGINAL
EPA SAMPLE NO.

CKY35

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 18347 SAS No.:

SDG No.: CKY28

Matrix: (soil/water) SOIL

Lab Sample ID: 499468

Sample wt/vol: 30.00(g/ml)G

Lab File ID:

% Moisture: 3 decanted: (Y/N) N

Date Received: 06/24/92

Extraction: (SepF/Cont/Sonc) SONC

Date Extracted: 06/26/92

Concentrated Extract Volume: 5000(uL)

Date Analyzed: 07/07/92

Injection Volume: 2.0(uL)

Dilution Factor: 1

GPC Cleanup: (Y/N) Y pH: 6.9

Sulfur Cleanup: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/KG</u>	Q
319-84-6-----	alpha-BHC	0.100	J
319-85-7-----	beta-BHC	1.8	U
319-86-8-----	delta-BHC	1.8	U
58-89-9-----	gamma-BHC (Lindane)	0.41	JP
76-44-8-----	Heptachlor	0.050	JP
309-00-2-----	Aldrin	1.8	U
1024-57-3-----	Heptachlor epoxide	0.10	JP
959-98-8-----	Endosulfan I	0.30	JP
60-57-1-----	Dieldrin	3.4	U
72-55-9-----	4,4'-DDE	3.4	U
72-20-8-----	Endrin	0.27	JP
33213-65-9-----	Endosulfan II	0.22	JP
72-54-8-----	4,4'-DDD	3.4	U
1031-07-8-----	Endosulfan sulfate	3.4	U
50-29-3-----	4,4'-DDT	3.4	U
72-43-5-----	Methoxychlor	9.4	JPB
53494-70-5-----	Endrin ketone	0.47	J
7421-93-4-----	Endrin aldehyde	3.4	U
5103-71-9-----	alpha-Chlordane	0.39	JP
5103-74-2-----	gamma-Chlordane	0.41	J
8001-35-2-----	Toxaphene	180	U
12674-11-2-----	Aroclor-1016	34	U
11104-28-2-----	Aroclor-1221	69	U
11141-16-5-----	Aroclor-1232	34	U
53469-21-9-----	Aroclor-1242	34	U
12672-29-6-----	Aroclor-1248	34	U
11097-69-1-----	Aroclor-1254	34	U
11096-82-5-----	Aroclor-1260	34	U

ID
PESTICIDE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

(Reg)

CKY36

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 18347 SAS No.:

SDG No.: CKY28

Matrix: (soil/water) SOIL

Lab Sample ID: 499470

Sample wt/vol: 30.00(g/ml)G

Lab File ID:

% Moisture: 10 decanted: (Y/N) N

Date Received: 06/24/92

Extraction: (SepF/Cont/Sonc) SONC

Date Extracted: 06/26/92

Concentrated Extract Volume: 5000(uL)

Date Analyzed: 07/07/92

Injection Volume: 2.0(uL)

Dilution Factor: 1

GPC Cleanup: (Y/N) Y pH: 7.0

Sulfur Cleanup: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) UG/KG	Q
319-84-6-----	alpha-BHC	1.9	U
319-85-7-----	beta-BHC	1.9	U
319-86-8-----	delta-BHC	1.9	U
58-89-9-----	gamma-BHC (Lindane)	0.37	JP
76-44-8-----	Heptachlor	1.9	U
309-00-2-----	Aldrin	1.9	U
1024-57-3-----	Heptachlor epoxide	1.9	U
959-98-8-----	Endosulfan I	1.9	U
60-57-1-----	Dieldrin	3.7	U
72-55-9-----	4,4'-DDE	0.97	J
72-20-8-----	Endrin	0.51	JP
33213-65-9-----	Endosulfan II	3.7	U
72-54-8-----	4,4'-DDD	3.7	U
1031-07-8-----	Endosulfan sulfate	3.7	U
50-29-3-----	4,4'-DDT	1.2	JP
72-43-5-----	Methoxychlor	0.99	JPB
53494-70-5-----	Endrin ketone	0.17	JP
7421-93-4-----	Endrin aldehyde	3.7	U
5103-71-9-----	alpha-Chlordane	1.9	U
5103-74-2-----	gamma-Chlordane	0.45	JP
8001-35-2-----	Toxaphene	190	U
12674-11-2-----	Aroclor-1016	37	U
11104-28-2-----	Aroclor-1221	74	U
11141-16-5-----	Aroclor-1232	37	U
53469-21-9-----	Aroclor-1242	37	U
12672-29-6-----	Aroclor-1248	37	U
11097-69-1-----	Aroclor-1254	37	U
11096-82-5-----	Aroclor-1260	37	U

1D
PESTICIDE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

02/
Recd

CKY37

Lab Name: COMPUCHEM, RTPContract: 68D10083Lab Code: COMPU Case No.: 18347 SAS No.:SDG No.: CKY28Matrix: (soil/water) SOILLab Sample ID: 499471Sample wt/vol: 30.00(g/ml)G

Lab File ID:

% Moisture: 8 decanted: (Y/N) NDate Received: 06/24/92Extraction: (SepF/Cont/Sonc) SONCDate Extracted: 06/26/92Concentrated Extract Volume: 5000(uL)Date Analyzed: 07/07/92Injection Volume: 2.0(uL)Dilution Factor: 1GPC Cleanup: (Y/N) Y pH: 6.1Sulfur Cleanup: (Y/N) N

CONCENTRATION UNITS:

CAS NO. COMPOUND (ug/L or ug/Kg) UG/KG Q

319-84-6-----	alpha-BHC	1.8	U
319-85-7-----	beta-BHC	1.8	U
319-86-8-----	delta-BHC	1.8	U
58-89-9-----	gamma-BHC (Lindane)	0.12	JP
76-44-8-----	Heptachlor	1.8	U
309-00-2-----	Aldrin	1.8	U
1024-57-3-----	Heptachlor epoxide	1.8	U
959-98-8-----	Endosulfan I	1.8	U
60-57-1-----	Dieldrin	3.6	U
72-55-9-----	4,4'-DDE	3.6	U
72-20-8-----	Endrin	3.6	U
33213-65-9-----	Endosulfan II	3.6	U
72-54-8-----	4,4'-DDD	3.6	U
1031-07-8-----	Endosulfan sulfate	3.6	U
50-29-3-----	4,4'-DDT	3.6	U
72-43-5-----	Methoxychlor	18	U
53494-70-5-----	Endrin ketone	0.12	J
7421-93-4-----	Endrin aldehyde	3.6	U
5103-71-9-----	alpha-Chlordane	1.8	U
5103-74-2-----	gamma-Chlordane	1.8	U
8001-35-2-----	Toxaphene	180	U
12674-11-2-----	Aroclor-1016	36	U
11104-28-2-----	Aroclor-1221	73	U
11141-16-5-----	Aroclor-1232	36	U
53469-21-9-----	Aroclor-1242	36	U
12672-29-6-----	Aroclor-1248	36	U
11097-69-1-----	Aroclor-1254	36	U
11096-82-5-----	Aroclor-1260	36	U

ID
PESTICIDE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CKY38

Lab Name: COMPUCHEM, RTPContract: 68D10083Lab Code: COMPU Case No.: 18347 SAS No.:SDG No.: CKY28Matrix: (soil/water) SOILLab Sample ID: 499473Sample wt/vol: 30.00(g/ml)G

Lab File ID:

% Moisture: 12 decanted: (Y/N) NDate Received: 06/24/92Extraction: (SepF/Cont/Sonc) SONCDate Extracted: 06/26/92Concentrated Extract Volume: 5000(uL)Date Analyzed: 07/07/92Injection Volume: 2.0(uL)Dilution Factor: 1GPC Cleanup: (Y/N) Y pH: 6.7Sulfur Cleanup: (Y/N) NCONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG

Q

<u>CAS NO.</u>	<u>COMPOUND</u>		
<u>319-84-6</u>	<u>alpha-BHC</u>	<u>1.9</u>	<u>U</u>
<u>319-85-7</u>	<u>beta-BHC</u>	<u>1.9</u>	<u>U</u>
<u>319-86-8</u>	<u>delta-BHC</u>	<u>1.9</u>	<u>U</u>
<u>58-89-9</u>	<u>gamma-BHC (Lindane)</u>	<u>1.9</u>	<u>U</u>
<u>76-44-8</u>	<u>Heptachlor</u>	<u>0.094</u>	<u>JP</u>
<u>309-00-2</u>	<u>Aldrin</u>	<u>0.37</u>	<u>JP</u>
<u>1024-57-3</u>	<u>Heptachlor epoxide</u>	<u>0.26</u>	<u>JP</u>
<u>959-98-8</u>	<u>Endosulfan I</u>	<u>0.34</u>	<u>J</u>
<u>60-57-1</u>	<u>Dieldrin</u>	<u>0.33</u>	<u>JP</u>
<u>72-55-9</u>	<u>4,4'-DDE</u>	<u>3.8</u>	<u>U</u>
<u>72-20-8</u>	<u>Endrin</u>	<u>0.56</u>	<u>JP</u>
<u>33213-65-9</u>	<u>Endosulfan II</u>	<u>3.8</u>	<u>U</u>
<u>72-54-8</u>	<u>4,4'-DDD</u>	<u>1.3</u>	<u>JP</u>
<u>1031-07-8</u>	<u>Endosulfan sulfate</u>	<u>3.8</u>	<u>U</u>
<u>50-29-3</u>	<u>4,4'-DDT</u>	<u>3.8</u>	<u>U</u>
<u>72-43-5</u>	<u>Methoxychlor</u>	<u>0.67</u>	<u>JPB</u>
<u>53494-70-5</u>	<u>Endrin ketone</u>	<u>3.8</u>	<u>U</u>
<u>7421-93-4</u>	<u>Endrin aldehyde</u>	<u>3.8</u>	<u>U</u>
<u>5103-71-9</u>	<u>alpha-Chlordane</u>	<u>0.48</u>	<u>JP</u>
<u>5103-74-2</u>	<u>gamma-Chlordane</u>	<u>0.51</u>	<u>JP</u>
<u>8001-35-2</u>	<u>Toxaphene</u>	<u>190</u>	<u>U</u>
<u>12674-11-2</u>	<u>Aroclor-1016</u>	<u>38</u>	<u>U</u>
<u>11104-28-2</u>	<u>Aroclor-1221</u>	<u>76</u>	<u>U</u>
<u>11141-16-5</u>	<u>Aroclor-1232</u>	<u>38</u>	<u>U</u>
<u>53469-21-9</u>	<u>Aroclor-1242</u>	<u>38</u>	<u>U</u>
<u>12672-29-6</u>	<u>Aroclor-1248</u>	<u>38</u>	<u>U</u>
<u>11097-69-1</u>	<u>Aroclor-1254</u>	<u>38</u>	<u>U</u>
<u>11096-82-5</u>	<u>Aroclor-1260</u>	<u>38</u>	<u>U</u>

ORIGINAL
(Red)

Appendix D

Reviewed and Corrected Tentatively Identified Compounds

1E
VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

OP
SAMPLE
REC'D

CKY14

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 18347

SAS No.: _____ SDG No.: CKY14

Matrix: (soil/water) WATER

Lab Sample ID: 499425

Sample wt/vol: 5.0 (g/mL) ML

Lab File ID: CN099425A56

Level: (low/med) LOW

Date Received: 06/24/92

% Moisture: not dec. _____

Date Analyzed: 06/25/92

Column: DB624 ID: 0.530 (mm)

Dilution Factor: 1.0

Soil Extract Volume: _____ (uL)

Soil Aliquot Volume: _____ (uL)

Number TICs found: 0

CONCENTRATION UNITS:

(ug/L or ug/Kg) UG/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
=====	=====	=====	=====	=====

1E
VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

CKY15

Lab Name: COMPUCHEM, RTP Contract: 68D10083

Lab Code: COMPU Case No.: 18347 SAS No.: _____ SDG No.: CKY14

Matrix: (soil/water) WATER

Lab Sample ID: 499435

Sample wt/vol: 5.0 (g/mL) ML

Lab File ID: CN099435B56

Level: (low/med) LOW

Date Received: 06/24/92

% Moisture: not dec. _____

Date Analyzed: 06/24/92

GC Column: DB624 ID: 0.530 (mm)

Dilution Factor: 1.0

Soil Extract Volume: _____ (uL)

Soil Aliquot Volume: _____ (uL)

Number TICs found: 0

CONCENTRATION UNITS:

(ug/L or ug/Kg) UG/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
=====	=====	=====	=====	=====

1E
VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

ORIGINIAL
EPA SAMPLE NO.

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

CKY16

Lab Code: COMPU Case No.: 18347 SAS No.: _____ SDG No.: CKY14

Matrix: (soil/water) WATER

Lab Sample ID: 499436

Sample wt/vol: 5.0 (g/mL) ML

Lab File ID: CN099436B56

Level: (low/med) LOW

Date Received: 06/24/92

% Moisture: not dec.

Date Analyzed: 06/24/92

GC Column: DB624 ID: 0.530 (mm)

Dilution Factor: 1.0

Soil Extract Volume: _____ (uL)

Soil Aliquot Volume: _____ (uL)

Number TICs found: 0

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
=====	=====	=====	=====	=====

VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

CKY17

Lab Name: COMPUCHEM RTPContract: 68D10083Lab Code: COMPU Case No.: 18347 SAS No.: _____ SDG No.: CKY14Matrix: (soil/water) WATERLab Sample ID: 499437Sample wt/vol: 5.0 (g/mL) MLLab File ID: CN099437A56Level: (low/med) LOWDate Received: 06/24/92

% Moisture: not dec. _____

Date Analyzed: 06/25/92GC Column: DB624 ID: 0.530 (mm)Dilution Factor: 1.0

Oil Extract Volume: _____ (uL)

Soil Aliquot Volume: _____ (uL)

Number TICs found: 0

CONCENTRATION UNITS:

(ug/L or ug/Kg) UG/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
=====	=====	=====	=====	=====

1E
VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

ORIGINAL
EPA SAMPLE NO.

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

CKY18

Lab Code: COMPU Case No.: 18347

SAS No.: _____ SDG No.: CKY14

Matrix: (soil/water) WATER

Lab Sample ID: 499439

Sample wt/vol: 5.0 (g/mL) ML

Lab File ID: CR099439C56

Level: (low/med) LOW

Date Received: 06/24/92

% Moisture: not dec. _____

Date Analyzed: 06/26/92

Column: DB624 ID: 0.530 (mm)

Dilution Factor: 1.0

Soil Extract Volume: _____ (uL)

Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

(ug/L or ug/Kg) UG/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
-----	-----	-----	-----	-----

1E
VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.
(Med)

CKY19

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 18347

SAS No.: _____ SDG No.: CKY14

Matrix: (soil/water) WATER

Lab Sample ID: 499440

Sample wt/vol: 5.0 (g/mL) ML

Lab File ID: CN099440A56

Level: (low/med) LOW

Date Received: 06/24/92

% Moisture: not dec.

Date Analyzed: 06/25/92

GC Column: DB624 ID: 0.530 (mm)

Dilution Factor: 1.0

Soil Extract Volume: _____ (uL)

Soil Aliquot Volume: _____ (uL)

Number TICs found: 0

CONCENTRATION UNITS:

(ug/L or ug/Kg) UG/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
=====	=====	=====	=====	=====

1E

EPA SAMPLE NO.

VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

Lab Name: COMPUCHEM, RTPContract: 68D10083

CKY20

Lab Code: COMPU Case No.: 18347

SAS No.: _____

SDG No.: CKY14Matrix: (soil/water) WATERLab Sample ID: 499441Sample wt/vol: 5.0 (g/mL) MLLab File ID: CN099441A56Level: (low/med) LOWDate Received: 06/24/92

Moisture: not dec. _____

Date Analyzed: 06/25/92Column: DB624 ID: 0.530 (mm)Dilution Factor: 1.0

Soil Extract Volume: _____ (uL)

Soil Aliquot Volume: _____ (uL)

Number TICs found: 0

CONCENTRATION UNITS:

(ug/L or ug/Kg) UG/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
=====	=====	=====	=====	=====

1E
VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.
(MSD)

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

CKY22

Lab Code: COMPU Case No.: 18347

SAS No.: _____ SDG No.: CKY14

Matrix: (soil/water) WATER

Lab Sample ID: 499442

Sample wt/vol: 5.0 (g/mL) ML

Lab File ID: CN099442A56

Level: (low/med) LOW

Date Received: 06/24/92

% Moisture: not dec. _____

Date Analyzed: 06/25/92

GC Column: DB624 ID: 0.530 (mm)

Dilution Factor: 1.0

Soil Extract Volume: _____ (uL)

Soil Aliquot Volume: _____ (uL)

Number TICs found: 0

CONCENTRATION UNITS:

(ug/L or ug/Kg) UG/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
=====	=====	=====	=====	=====

1E
VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

CKY23

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 18347

SAS No.: _____ SDG No.: CKY14

Matrix: (soil/water) WATER

Lab Sample ID: 499444

Sample wt/vol: 5.0 (g/mL) ML

Lab File ID: CN099444A56

Level: (low/med) LOW

Date Received: 06/24/92

% Moisture: not dec.

Date Analyzed: 06/25/92

GC Column: DB624 ID: 0.530 (mm)

Dilution Factor: 1.0

Soil Extract Volume: _____ (uL)

Soil Aliquot Volume: _____ (uL)

Number TICs found: 0

CONCENTRATION UNITS:

(ug/L or ug/Kg) UG/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
=====	=====	=====	=====	=====

1E
VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

ORIGINAL

(Red)

CKY24

Lab Name: COMPUCHEM RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 18347 SAS No.: _____ SDG No.: CKY14

Matrix: (soil/water) WATER

Lab Sample ID: 499445

Sample wt/vol: 5.0 (g/mL) ML

Lab File ID: CN099445A56

Level: (low/med) LOW

Date Received: 06/24/92

% Moisture: not dec. _____

Date Analyzed: 06/25/92

Column: DB624 ID: 0.530 (mm)

Dilution Factor: 1.0

Soil Extract Volume: _____ (uL)

Soil Aliquot Volume: _____ (uL)

Number TICs found: 0

CONCENTRATION UNITS:

(ug/L or ug/Kg) UG/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
-----	-----	-----	-----	-----

1E
VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

ORIGINAL

(Rev)

CKY25

Lab Name: COMPUCHEM, RTP Contract: 68D10083

Lab Code: COMPU Case No.: 18347 SAS No.: _____ SDG No.: CKY14

Matrix: (soil/water) WATER

Lab Sample ID: 499447

Sample wt/vol: 5.0 (g/mL) ML

Lab File ID: CN099447B56

Level: (low/med) LOW

Date Received: 06/24/92

% Moisture: not dec. _____

Date Analyzed: 06/25/92

GC Column: DB624 ID: 0.530 (mm)

Dilution Factor: 1.0

Soil Extract Volume: _____ (uL)

Soil Aliquot Volume: _____ (uL)

Number TICs found: 0

CONCENTRATION UNITS:

(ug/L or ug/Kg) UG/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
=====	=====	=====	=====	=====

1E
VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA/SAMPLE NO.

(Reg)

CKY26

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 18347

SAS No.: _____ SDG No.: CKY14

Matrix: (soil/water) WATER

Lab Sample ID: 499449

Sample wt/vol: 5.0 (g/mL) ML

Lab File ID: CR099449C56

Level: (low/med) LOW

Date Received: 06/24/92

% Moisture: not dec. _____

Date Analyzed: 06/26/92

Column: DB624 ID: 0.530 (mm)

Dilution Factor: 1.0

Soil Extract Volume: _____ (uL)

Soil Aliquot Volume: _____ (uL)

Number TICs found: 0

CONCENTRATION UNITS:

(ug/L or ug/Kg) UG/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
=====	=====	=====	=====	=====

1E
VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

(Rev. 1)

CKY28

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 18347

SAS No.: _____ SDG No.: CKY28

Matrix: (soil/water) SOIL

Lab Sample ID: 499452

Sample wt/vol: 5.0 (g/mL) G

Lab File ID: GH099452B54

Level: (low/med) LOW

Date Received: 06/24/92

% Moisture: not dec. 19

Date Analyzed: 06/24/92

GC Column: DB-624 ID: 0.530 (mm)

Dilution Factor: 1.0

Soil Extract Volume: _____ (uL)

Soil Aliquot Volume: _____ (uL)

Number TICs found: 1

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1.	LABORATORY ARTIFACT	2.20	36	J

1E
VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

ORIGINAL

(Red)

CKY29

Lab Name: COMPUCHEM, RTP

Contract: 58D10083

Lab Code: COMPU Case No.: 18347 SAS No.: _____ SDG No.: CKY28

Matrix: (soil/water) SOIL

Lab Sample ID: 499459

Sample wt/vol: 5.0 (g/mL) G

Lab File ID: GH099459B54

Level: (low/med) LOW

Date Received: 06/24/92

% Moisture: not dec. 17

Date Analyzed: 06/24/92

GC Column: DB-624 ID: 0.530 (mm)

Dilution Factor: 1.0

Soil Extract Volume: _____ (uL)

Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

(ug/L or ug/Kg) UG/KG

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1.	LABORATORY ARTIFACT	1.55	200	J
2.	LABORATORY ARTIFACT	2.18	36	J

1E
VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

FINAL

CKY30

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 18347

SAS No.: _____ SDG No.: CKY28

Matrix: (soil/water) SOIL

Lab Sample ID: 499463

Sample wt/vol: 5.0 (g/mL) G

Lab File ID: GH099463B54

Level: (low/med) LOW

Date Received: 06/24/92

% Moisture: not dec. 40

Date Analyzed: 06/24/92

Column: DB-624 ID: 0.530 (mm)

Dilution Factor: 1.0

Soil Extract Volume: _____ (uL)

Soil Aliquot Volume: _____ (uL)

Number TICs found: 1

CONCENTRATION UNITS:

(ug/L or ug/Kg) UG/KG

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1.	LABORATORY ARTIFACT	2.17	130	BJ

1E
 VOLATILE ORGANICS ANALYSIS DATA SHEET
 TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

ORIGINAL

(Red)

CKY31

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 18347

SAS No.: _____ SDG No.: CKY28

Matrix: (soil/water) SOIL

Lab Sample ID: 499464

Sample wt/vol: 5.0 (g/mL) G

Lab File ID: GH099464B54

Level: (low/med) LOW

Date Received: 06/24/92

% Moisture: not dec. 23

Date Analyzed: 06/24/92

Column: DB-624 ID: 0.530 (mm)

Dilution Factor: 1.0

Soil Extract Volume: _____ (uL)

Soil Aliquot Volume: _____ (uL)

Number TICs found: 1

CONCENTRATION UNITS:

(ug/L or ug/Kg) UG/KG

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
-1.	LABORATORY-ARTIFACT	2.20	38	J

1E

EPA SAMPLE NO.

VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

Lab Name: COMPUCHEM, RTPContract: 68D10083

CKY32

Lab Code: COMPU Case No.: 18347SAS No.: _____ SDG No.: CKY28Matrix: (soil/water) SOILLab Sample ID: 499465Sample wt/vol: 5.0 (g/mL) GLab File ID: GR099465A54Level: (low/med) LOWDate Received: 06/24/92% Moisture: not dec. 18Date Analyzed: 06/25/92Column: DB-624 ID: 0.530 (mm)Dilution Factor: 1.0

Soil Extract Volume: _____ (uL)

Soil Aliquot Volume: _____ (uL)

Number TICs found: 1

CONCENTRATION UNITS:

(ug/L or ug/Kg) UG/KG

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
-1.	UNKNOWN	2.17	43	J K C

1E
 VOLATILE ORGANICS ANALYSIS DATA SHEET
 TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

ORIGINAL
CKY28

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 18347

SAS No.: _____ SDG No.: CKY28

Matrix: (soil/water) SOIL

Lab Sample ID: 499466

Sample wt/vol: 5.0 (g/mL) G

Lab File ID: GR099466A54

Level: (low/med) LOW

Date Received: 06/24/92

% Moisture: not dec. 11

Date Analyzed: 06/25/92

GC Column: DB-624 ID: 0.530 (mm)

Dilution Factor: 1.0

Soil Extract Volume: _____ (uL)

Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

(ug/L or ug/Kg) UG/KG

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1.	LABORATORY ARTIFACT	2.18	11	BJ

1E
VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

(Randy)

CKY34

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU

Case No.: 18347

SAS No.: _____

SDG No.: CKY28

Matrix: (soil/water) SOIL

Lab Sample ID: 499467

Sample wt/vol: 5.0 (g/mL) G

Lab File ID: GH099467B54

Level: (low/med) LOW

Date Received: 06/24/92

% Moisture: not dec. 13

Date Analyzed: 06/24/92

GC Column: DB-624 ID: 0.530 (mm)

Dilution Factor: 1.0

Soil Extract Volume: _____ (uL)

Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

(ug/L or ug/Kg) UG/KG

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1.	LABORATORY ARTIFACT	2.20	29	J H

VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

ORIGINAL

10-01

CKY35

Lab Name: COMPUCHEM, RTPContract: 68D10083Lab Code: COMPU Case No.: 18347SAS No.: _____ SDG No.: CKY28Matrix: (soil/water) SOILLab Sample ID: 499468Sample wt/vol: 5.0 (g/mL) GLab File ID: GH099468B54Level: (low/med) LOWDate Received: 06/24/92% Moisture: not dec. 3Date Analyzed: 06/25/92GC Column: DB-624 ID: 0.530 (mm)Dilution Factor: 1.0

Soil Extract Volume: _____ (uL)

Soil Aliquot Volume: _____ (uL)

Number TICs found: 1CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
-1.	LABORATORY ARTIFACT	2.20	34	J

1E
VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

ORIGINAL
COPY

CKY36

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 18347 SAS No.: _____ SDG No.: CKY28

Matrix: (soil/water) SOIL

Lab Sample ID: 499470

Sample wt/vol: 5.0 (g/mL) G

Lab File ID: GR099470A54

Level: (low/med) LOW

Date Received: 06/24/92

% Moisture: not dec. 10

Date Analyzed: 06/25/92

GC Column: DB-624 ID: 0.530 (mm)

Dilution Factor: 1.0

Soil Extract Volume: _____ (uL)

Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

(ug/L or ug/Kg) UG/KG

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1.	LABORATORY ARTIFACT	1.55	43	J }
2.	LABORATORY ARTIFACT	2.18	14	J }

1E
VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

CKY37

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 18347 SAS No.: _____ SDG No.: CKY28

Matrix: (soil/water) SOIL

Lab Sample ID: 499471

Sample wt/vol: 5.0 (g/mL) G

Lab File ID: GH099471B54

Level: (low/med) LOW

Date Received: 06/24/92

% Moisture: not dec. 8

Date Analyzed: 06/25/92

Column: DB-624 ID: 0.530 (mm)

Dilution Factor: 1.0

Soil Extract Volume: _____ (uL)

Soil Aliquot Volume: _____ (uL)

Number TICs found: 1

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1.	LABORATORY-ARTIFACT	2.18	30	J

1E
VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

(Pacal)

Lab Name: COMPUCHEM RTP

Contract: 68D10083

CKY38

Lab Code: COMPU Case No.: 18347

SAS No.: _____ SDG No.: CKY28

Matrix: (soil/water) SOIL

Lab Sample ID: 499473

Sample wt/vol: 5.0 (g/mL) G

Lab File ID: GH099473C54

Level: (low/med) LOW

Date Received: 06/24/92

% Moisture: not dec. 12

Date Analyzed: 06/25/92

Column: DB-624 ID: 0.530 (mm)

Dilution Factor: 1.0

Soil Extract Volume: _____ (uL)

Soil Aliquot Volume: _____ (uL)

Number TICs found: 1

CONCENTRATION UNITS:

(ug/L or ug/Kg) UG/KG

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1.	LABORATORY ARTIFACT	2.15	54	BJ 91

1F
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA 04 SAMPLE NO.

(Reg)

CKY14

Lab Name: COMPUCHEM, RTP Contract: 68D10083

Lab Code: COMPU Case No.: 18347 SAS No.: _____ SDG No.: CKY14

Matrix: (soil/water) WATER Lab Sample ID: 499425

Sample wt/vol: 1000 (g/mL) ML Lab File ID: GH099425B02

Level: (low/med) LOW Date Received: 06/24/92

% Moisture: _____ decanted: (Y/N) _____ Date Extracted: 06/26/92

Concentrated Extract Volume: 1000 (uL) Date Analyzed: 06/29/92

Injection Volume: 2.0 (uL) Dilution Factor: 1.0

GPC Cleanup: (Y/N) N pH: _____

Number TICs found: 7

CONCENTRATION UNITS:

(ug/L or ug/Kg) UG/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1.	BLANK CONTAMINANT	4.80	26	BJ } H ₂ O
2.	BLANK CONTAMINANT	5.13	11	BJ } 91 mfc
3.	UNKNOWN	5.18	2	J
4.	UNKNOWN	10.80	11	J
5.	UNKNOWN	10.87	5	J
6.	LABORATORY ARTIFACT	14.07	7	J } 145
7.	LABORATORY ARTIFACT	15.37	17	BJ } 91 mfc

1F
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

ORIG.
EPA SAMPLE NO.
(Red)

CKY15

Lab Name: COMPUCHEM, RTP Contract: 68D10083

Lab Code: COMPU Case No.: 18347 SAS No.: _____ SDG No.: CKY14

Matrix: (soil/water) WATER Lab Sample ID: 499435

Sample wt/vol: 1000 (g/mL) ML Lab File ID: GH099435B02

Level: (low/med) LOW Date Received: 06/24/92

Moisture: _____ decanted: (Y/N) _____ Date Extracted: 06/26/92

Concentrated Extract Volume: 1000 (uL) Date Analyzed: 06/29/92

Injection Volume: 2.0 (uL) Dilution Factor: 1.0

GPC Cleanup: (Y/N) N pH: _____

CONCENTRATION UNITS:
Number TICs found: 5 (ug/L or ug/Kg) UG/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1.	UNKNOWN	4.42	6	J
2.	BLANK CONTAMINANT	4.80	17	BJ
3.	UNKNOWN	4.95	6	J
4.	UNKNOWN	5.07	4	J
5.	BLANK CONTAMINANT	5.13	14	BJ

1F
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.
(Red)

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

CKY16

Lab Code: COMPU Case No.: 18347

SAS No.: _____

SDG No.: CKY14

Matrix: (soil/water) WATER

Lab Sample ID: 499436

Sample wt/vol: 1000 (g/mL) ML

Lab File ID: GH099436B02

Level: (low/med) LOW

Date Received: 06/24/92

% Moisture: _____ decanted: (Y/N) _____

Date Extracted: 06/26/92

Concentrated Extract Volume: 1000 (uL)

Date Analyzed: 06/29/92

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) N pH: _____

Number TICs found: 6

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1.	UNKNOWN	4.48	4	J
2.	BLANK CONTAMINANT	4.67	13	BJ
3.	UNKNOWN	5.07	5	J
4.	LABORATORY ARTIFACT	14.10	6	J
5.	LABORATORY ARTIFACT	15.40	12	BJ
6.	UNKNOWN	16.65	5	J

1F
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.
(Regd.)

CKY17

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 18347

SAS No.: _____ SDG No.: CKY14

Matrix: (soil/water) WATER

Lab Sample ID: 499437

Sample wt/vol: 1000 (g/mL) ML

Lab File ID: GH099437B02

Level: (low/med) LOW

Date Received: 06/24/92

% Moisture: _____ decanted: (Y/N) _____

Date Extracted: 06/26/92

Concentrated Extract Volume: 1000 (uL)

Date Analyzed: 06/30/92

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) N pH: _____

CONCENTRATION UNITS:

(ug/L or ug/Kg) UG/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
-1.	LABORATORY-ARTIFACT	15.39	7	BJ

1F
SEMICVOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

ORIGIN
EPA SAMPLE NO.
(red)

CKY18

Lab Name: COMPUCHEM, RTP Contract: 68D10083

Lab Code: COMPU Case No.: 18347 SAS No.: _____ SDG No.: CKY14

Matrix: (soil/water) WATER Lab Sample ID: 499439

Sample wt/vol: 1000 (g/mL) ML Lab File ID: GH099439B02

Level: (low/med) LOW Date Received: 06/24/92

% Moisture: _____ decanted: (Y/N) _____ Date Extracted: 06/26/92

Concentrated Extract Volume: 1000 (uL) Date Analyzed: 06/29/92

Injection Volume: 2.0 (uL) Dilution Factor: 1.0

GPC Cleanup: (Y/N) N pH: _____

CONCENTRATION UNITS:
Number TICs found: 2 (ug/L or ug/Kg) UG/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1.	LABORATORY ARTIFACT	14.10	3	J-H
2.	LABORATORY ARTIFACT	15.42	27	BJ

1F
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

CKY19

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 18347

SAS No.: _____ SDG No.: CKY14

Matrix: (soil/water) WATER

Lab Sample ID: 499440

Sample wt/vol: 1000 (g/mL) ML

Lab File ID: GH099440B02

Level: (low/med) LOW

Date Received: 06/24/92

% Moisture: _____ decanted: (Y/N)

Date Extracted: 06/26/92

Concentrated Extract Volume: 1000 (uL)

Date Analyzed: 06/30/92

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) N pH: _____

CONCENTRATION UNITS:

Number TICs found: 4

(ug/L or ug/Kg) UG/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1.	UNKNOWN	4.50	4	J
2.	UNKNOWN	4.68	10	J
3.	UNKNOWN	5.08	4	J
4.	LABORATORY ARTIFACT	15.40	18	BJ

1F
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

^{09/04/92}
^(Med)
EPA SAMPLE NO.

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

CKY20

Lab Code: COMPU

Case No.: 18347

SAS No.: _____

SDG No.: CKY14

Matrix: (soil/water) WATER

Lab Sample ID: 499441

Sample wt/vol: 1000 (g/mL) ML

Lab File ID: GH099441B02

Level: (low/med) LOW

Date Received: 06/24/92

% Moisture: _____ decanted: (Y/N)

Date Extracted: 06/26/92

Concentrated Extract Volume: 1000 (uL)

Date Analyzed: 06/30/92

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) N pH: _____

Number TICs found: 17

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1.	UNKNOWN	5.10	9	J
2.	TRICHLOROPROPENE	5.18	3	J
3.	UNKNOWN CARBOXYLIC ACID	11.30	3	J
4.	UNKNOWN CARBOXYLIC ACID	12.45	20	J
5.	UNKNOWN	13.42	4	J
6.	UNKNOWN CARBOXYLIC ACID	13.49	9	J
7.	LAB. ARTIFACT	14.12	8	J
8.	LAB. ARTIFACT	15.42	22	BJ
9.	UNKNOWN	16.34	7	J
10.	UNKNOWN	16.67	7	J
11.	UNKNOWN	16.82	7	J
12.	UNKNOWN	17.35	5	J
13.	UNKNOWN	17.49	4	J
14.	UNKNOWN	19.15	68	J
15.	UNKNOWN	19.54	160	J
16.	UNKNOWN	20.02	9	J
17.	UNKNOWN	20.99	5	J

1F
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

ORIGINAL
EPA SAMPLE NO.

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

CKY22

Lab Code: COMPU Case No.: 18347 SAS No.: _____ SDG No.: CKY14

Matrix: (soil/water) WATER

Lab Sample ID: 499442

Sample wt/vol: 1000 (g/mL) ML

Lab File ID: GH099442B02

Level: (low/med) LOW

Date Received: 06/24/92

% Moisture: _____ decanted: (Y/N) _____

Date Extracted: 06/26/92

Concentrated Extract Volume: 1000 (uL)

Date Analyzed: 06/30/92

Injection Volume: 2.0(uL) pH: _____

Dilution Factor: 1.0

GPC Cleanup: (Y/N) N

Number TICs found: 3

CONCENTRATION UNITS:

(ug/L or ug/Kg) UG/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1.	BLANK CONTAMINANT	4.70	10	BJ
2.	BLANK CONTAMINANT	5.08	6	BJ
3.	LABORATORY ARTIFACT	15.42	12	BJ

1F
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

*original
med*

CKY23

Lab Name: COMPUCHEM, RTP Contract: 68D10083

Lab Code: COMPU Case No.: 18347 SAS No.: _____ SDG No.: CKY14

Matrix: (soil/water) WATER Lab Sample ID: 499444

Sample wt/vol: 1000 (g/mL) ML Lab File ID: GH099444B02

Level: (low/med) LOW Date Received: 06/24/92

% Moisture: _____ decanted: (Y/N) _____ Date Extracted: 06/26/92

Concentrated Extract Volume: 1000 (uL) Date Analyzed: 06/30/92

Injection Volume: 2.0 (uL) Dilution Factor: 1.0

GPC Cleanup: (Y/N) N pH: _____

Number TICs found: 2

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1.	BLANK CONTAMINANT	5.05	3	BJ 14
2.	LABORATORY ARTIFACT	15.37	5	BJ 9

**SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS**

*ORIGINAL
(Red)*
CKY24

Lab Name: COMPUCHEM, RTP Contract: 68D10083

Lab Code: COMPU Case No.: 18347 SAS No.: _____ SDG No.: CKY14

Matrix: (soil/water) WATER Lab Sample ID: 499445

Sample wt/vol: 1000 (g/mL) ML Lab File ID: GH099445A07

Level: (low/med) LOW Date Received: 06/24/92

% Moisture: _____ decanted: (Y/N) _____ Date Extracted: 06/25/92

Concentrated Extract Volume: 1000 (uL) Date Analyzed: 06/30/92

Injection Volume: 2.0 (uL) Dilution Factor: 1.0

GPC Cleanup: (Y/N) N pH: _____

Number TICs found: 3 CONCENTRATION UNITS:

(ug/L or ug/Kg) UG/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1.	BLANK CONTAMINANT	4.65	2	BJ
2.	UNKNOWN	13.57	5	J
3.	UNKNOWN	13.67	25	J

1F
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

ORIGINAL

(Red)

CKY25

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 18347

SAS No.: _____

SDG No.: CKY14

Matrix: (soil/water) WATER

Lab Sample ID: 499447

Sample wt/vol: 1000 (g/mL) ML

Lab File ID: GH099447A07

Level: (low/med) LOW

Date Received: 06/24/92

% Moisture: _____ decanted: (Y/N) _____

Date Extracted: 06/25/92

Concentrated Extract Volume: 1000 (uL)

Date Analyzed: 06/30/92

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) N pH: _____

CONCENTRATION UNITS:

Number TICs found: 0

(ug/L or ug/Kg) UG/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
=====	=====	=====	=====	=====

1F
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

ORIGINAL

(Red)

CKY26

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 18347

SAS No.: _____ SDG No.: CKY14

Matrix: (soil/water) WATER

Lab Sample ID: 499449

Sample wt/vol: 1000 (g/mL) ML

Lab File ID: GH099449A07

Level: (low/med) LOW

Date Received: 06/24/92

% Moisture: _____ decanted: (Y/N) _____

Date Extracted: 06/25/92

Concentrated Extract Volume: 1000 (uL)

Date Analyzed: 06/30/92

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) N pH: _____

Number TICs found: 3

CONCENTRATION UNITS:

(ug/L or ug/Kg) UG/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1.	UNKNOWN	4.67	7	J
2.	UNKNOWN	13.39	7	J
3.	UNKNOWN	14.44	18	J

1F
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

FINAL
REGD

CKY28

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 18347

SAS No.: _____ SDG No.: CKY28

Matrix: (soil/water) SOIL

Lab Sample ID: 499452

Sample wt/vol: 30.1 (g/mL) G

Lab File ID: GR099452A57

Level: (low/med) LOW

Date Received: 06/24/92

% Moisture: 19 decanted: (Y/N) N

Date Extracted: 07/07/92

Concentrated Extract Volume: 500.0 (uL)

Date Analyzed: 07/09/92

Injection Volume: 2.0(uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 7.2

Number TICs found: 15

CONCENTRATION UNITS:

(ug/L or ug/Kg) UG/KG

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1.	UNKNOWN	6.40	120	J
2.	BLANK CONTAMINANT	6.48	160	BJ
3.	BLANK CONTAMINANT	6.55	200	BJ
4.	UNKNOWN ALKANE	8.47	370	J
5.	UNKNOWN	12.24	82	J
6.	UNKNOWN HYDROCARBON	15.14	82	J
7.	UNKNOWN HYDROCARBON	17.10	82	J
8.	LABORATORY ARTIFACT	17.62	290	J
9.	UNKNOWN HYDROCARBON	18.29	120	J
10.	UNKNOWN HYDROCARBON	19.84	160	J
11.	UNKNOWN	21.30	120	J
12.	UNKNOWN HYDROCARBON	21.97	290	J
13.	UNKNOWN	24.14	120	J
14.	UNKNOWN	24.15	160	J
15.	UNKNOWN HYDROCARBON	25.11	200	J

1F
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

ORIGINAL

(Red)

CKY29

Lab Name: COMPUCHEM, RTP Contract: 68D10083

Lab Code: COMPU Case No.: 18347 SAS No.: SDG No.: CKY28

Matrix: (soil/water) SOIL

Lab Sample ID: 499459

Sample wt/vol: 30.7 (g/mL) G

Lab File ID: G2R99459A05

Level: (low/med) LOW

Date Received: 06/24/92

% Moisture: 17 decanted: (Y/N) N

Date Extracted: 07/10/92

Concentrated Extract Volume: 500.0 (uL)

Date Analyzed: 07/15/92

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 6.6

CONCENTRATION UNITS:

Number TICs found: 17

(ug/L or ug/Kg) UG/KG

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1.	BLANK CONTAMINANT	4.75	200	BJ
2.	TETRACHLOROETHANE	4.82	120	J
3.	UNKNOWN HYDROCARBON	6.98	120	J
4.	UNKNOWN HYDROCARBON	7.63	79	J
5.	UNKNOWN	7.68	430	J
6.	UNKNOWN	8.70	510	J
7.	UNKNOWN HYDROCARBON	8.84	120	J
8.	UNKNOWN HYDROCARBON	9.37	160	J
9.	UNKNOWN	9.75	390	J
10.	UNKNOWN	9.80	200	J
11.	UNKNOWN HYDROCARBON	10.39	200	J
12.	UNKNOWN + HYDROCARBON	10.85	200	J
13.	LABORATORY ARTIFACT	11.40	310	BJ
14.	LABORATORY ARTIFACT	12.85	200	BJ
15.	LABORATORY ARTIFACT	13.07	8600	BJ
16.	UNKNOWN	16.67	240	J
17.	UNKNOWN PAH	17.19	160	J

1F
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

PARTIAL
(Rev)

CKY30

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU

Case No.: 18347

SAS No.: _____

SDG No.: CKY28

Matrix: (soil/water) SOIL

Lab Sample ID: 499463

Sample wt/vol: 30.0 (g/mL) G

Lab File ID: GH099463A15

Level: (low/med) LOW

Date Received: 06/24/92

% Moisture: 40 decanted: (Y/N) N

Date Extracted: 06/26/92

Concentrated Extract Volume: 500.0 (uL)

Date Analyzed: 07/01/92

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 5.5

CONCENTRATION UNITS:

Number TICs found: 26

(ug/L or ug/Kg) UG/KG

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1.	UNKNOWN	5.10	1200	J
2.	BLANK-CONTAMINANT	5.20	5600	BJ
3.	BLANK-CONTAMINANT	5.25	670	BJ
4.	BLANK-CONTAMINANT	5.32	720	BJ
5.	BLANK CONTAMINANT	5.35	1200	BJ
6.	UNKNOWN	5.45	1300	J
7.	UNKNOWN	5.62	1100	J
8.	UNKNOWN	5.77	1700	J
9.	UNKNOWN	5.80	1000	J
10.	UNKNOWN	5.83	780	J
11.	UNKNOWN	6.33	610	J
12.	UNKNOWN	10.72	720	J
13.	UNKNOWN	11.25	780	J
14.	UNKNOWN HYDROCARBON	12.87	670	J
15.	UNKNOWN	13.05	560	J
16.	LABORATORY ARTIFACT	13.32	780	J
17.	LABORATORY ARTIFACT	13.37	14000	J
18.	UNKNOWN	13.62	830	J
19.	UNKNOWN HYDROCARBON	14.00	2700	J
20.	UNKNOWN	14.27	1200	J
21.	UNKNOWN HYDROCARBON	15.50	4000	J
22.	UNKNOWN	16.97	2300	J
23.	UNKNOWN HYDROCARBON	17.70	17000	J
24.	UNKNOWN	21.07	4600	J
25.	UNKNOWN	21.17	7800	J
26.	UNKNOWN	27.16	6700	J

1F
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

*Original
Ready*
CKY31

Lab Name: COMPUCHEM, RTP Contract: 68D10083

Lab Code: COMPU Case No.: 18347 SAS No.: _____ SDG No.: CKY28

Matrix: (soil/water) SOIL

Lab Sample ID: 499464

Sample wt/vol: 30.0 (g/mL) G

Lab File ID: GR099464A57

Level: (low/med) LOW

Date Received: 06/24/92

% Moisture: 23 decanted: (Y/N) N

Date Extracted: 07/07/92

Concentrated Extract Volume: 500.0 (uL)

Date Analyzed: 07/09/92

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 7.1

CONCENTRATION UNITS:

Number TICs found: 22

(ug/L or ug/Kg) UG/KG

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1.	UNKNOWN	5.98	1100	J
2.	UNKNOWN	6.43	130	J
3.	BLANK CONTAMINANT	6.50	170	BJ
4.	BLANK CONTAMINANT	6.58	260	BJ
5.	UNKNOWN ALKANE	8.49	170	J
6.	UNKNOWN	16.89	130	J
7.	UNKNOWN HYDROCARBON	17.12	86	J
8.	LABORATORY ARTIFACT	17.65	130	J
9.	UNKNOWN HYDROCARBON	18.32	220	J
10.	UNKNOWN	18.95	220	J
11.	UNKNOWN	19.37	130	J
12.	UNKNOWN HYDROCARBON	19.85	390	J
13.	UNKNOWN	19.99	260	J
14.	UNKNOWN	21.32	350	J
15.	UNKNOWN HYDROCARBON	22.00	950	J
16.	UNKNOWN	22.20	260	J
17.	UNKNOWN	23.39	130	J
18.	UNKNOWN	24.19	560	J
19.	UNKNOWN HYDROCARBON	25.12	650	J
20.	UNKNOWN HYDROCARBON	25.14	610	J
21.	UNKNOWN	28.47	390	J
22.	UNKNOWN	28.49	350	J

1F
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

*High
med*

CKY32

Lab Name: COMPUCHEM, RTP Contract: 68D10083

Lab Code: COMPU Case No.: 18347 SAS No.: _____ SDG No.: CKY28

Matrix: (soil/water) SOIL Lab Sample ID: 499465

Sample wt/vol: 30.0 (g/mL) G Lab File ID: GH099465A15

Level: (low/med) LOW Date Received: 06/24/92

% Moisture: 18 decanted: (Y/N) N Date Extracted: 06/26/92

Concentrated Extract Volume: 500.0 (uL) Date Analyzed: 07/01/92

Injection Volume: 2.0 (uL) Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 6.5

CONCENTRATION UNITS:

Number TICs found: 26 (ug/L or ug/Kg) UG/KG

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1.	UNKNOWN	5.02	200	J
2.	UNKNOWN	5.08	1400	J
3.	BLANK CONTAMINANT	5.20	5700	BJ
4.	BLANK CONTAMINANT	5.23	610	BJ
5.	BLANK CONTAMINANT	5.33	1300	BJ
6.	UNKNOWN	5.38	160	J
7.	UNKNOWN	5.45	610	J
8.	UNKNOWN	5.60	1300	J
9.	BLANK CONTAMINANT	5.78	1900	BJ
10.	UNKNOWN	5.83	320	J
11.	UNKNOWN	6.02	160	J
12.	UNKNOWN	6.22	450	J
13.	UNKNOWN	6.33	320	J
14.	UNKNOWN	6.60	200	J
15.	UNKNOWN	6.90	160	J
16.	UNKNOWN HYDROCARBON	8.99	120	J
17.	UNKNOWN HYDROCARBON	9.50	160	J
18.	UNKNOWN	11.22	1100	J
19.	UNKNOWN	11.25	730	J
20.	UNKNOWN	12.10	370	J
21.	LABORATORY ARTIFACT	13.29	370	J
22.	LABORATORY ARTIFACT	13.34	6900	J
23.	UNKNOWN HYDROCARBON	13.97	650	J
24.	UNKNOWN HYDROCARBON	15.45	970	J
25.	UNKNOWN	16.92	1300	J
26.	UNKNOWN	17.70	4000	J

ORIGINAL
(Red)

EPA SAMPLE NO.

1F

SEMICVOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

CKY33

Lab Name: COMPUCHEM, RTP Contract: 68D10083Lab Code: COMPU Case No.: 18347 SAS No.: SDG No.: CKY28Matrix: (soil/water) SOILLab Sample ID: 499466Sample wt/vol: 30.0 (g/mL) GLab File ID: GH099466A15Level: (low/med) LOWDate Received: 06/24/92% Moisture: 11 decanted: (Y/N) NDate Extracted: 06/26/92Concentrated Extract Volume: 500.0 (uL)Date Analyzed: 07/02/92Injection Volume: 2.0(uL)Dilution Factor: 1.0GPC Cleanup: (Y/N) Y pH: 5.9

CONCENTRATION UNITS:

Number TICs found: 26(ug/L or ug/Kg) UG/KG

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1.	UNKNOWN	5.05	150	J
2.	UNKNOWN	5.10	1200	J
3.	BLANK-CONTAMINANT	5.20	3500	BJ
4.	BLANK-CONTAMINANT	5.25	300	BJ
5.	BLANK-CONTAMINANT	5.32	370	BJ
6.	BLANK-CONTAMINANT	5.35	450	BJ
7.	UNKNOWN	5.47	450	J
8.	UNKNOWN	5.62	790	J
9.	BLANK-CONTAMINANT	5.80	1500	BJ
10.	UNKNOWN	6.35	260	J
11.	UNKNOWN	6.62	190	J
12.	UNKNOWN HYDROCARBON	7.28	260	J
13.	UNKNOWN	9.75	300	J
14.	UNKNOWN	9.99	190	J
15.	UNKNOWN	10.14	150	J
16.	UNKNOWN	10.74	260	J
17.	UNKNOWN	11.27	560	J
18.	UNKNOWN HYDROCARBON	11.47	190	J
19.	UNKNOWN	11.94	370	J
20.	LABORATORY-ARTIFACT	13.39	7500	J
21.	UNKNOWN	13.64	260	J
22.	UNKNOWN	14.02	520	J
23.	UNKNOWN	14.29	340	J
24.	UNKNOWN	15.52	450	J
25.	UNKNOWN	17.72	2200	J
26.	UNKNOWN	17.84	2000	J

1F
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

ORIGINAL
(Rev)

CKY34

Lab Name: COMPUCHEM, RTP Contract: 68D10083

Lab Code: COMPU Case No.: 18347 SAS No.: _____ SDG No.: CKY28

Matrix: (soil/water) SOIL

Lab Sample ID: 499467

Sample wt/vol: 30.0 (g/mL) G

Lab File ID: GR099467B57

Level: (low/med) LOW

Date Received: 06/24/92

% Moisture: 13 decanted: (Y/N) N

Date Extracted: 07/07/92

Concentrated Extract Volume: 500.0 (uL)

Date Analyzed: 07/09/92

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 6.5

CONCENTRATION UNITS:

Number TICs found: 6

(ug/L or ug/Kg) UG/KG

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1.	UNKNOWN HYDROCARBON	19.82	150	J
2.	UNKNOWN	21.27	270	J
3.	UNKNOWN HYDROCARBON	21.97	500	J
4.	UNKNOWN	23.30	730	J
5.	UNKNOWN	24.12	150	J
6.	UNKNOWN HYDROCARBON	25.07	460	J

1F
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

CKY35

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU

Case No.: 18347

SAS No.: _____

SDG No.: CKY28

Matrix: (soil/water) SOIL

Lab Sample ID: 499468

Sample wt/vol: 30.0 (g/mL) G

Lab File ID: GR099468A57

Level: (low/med) LOW

Date Received: 06/24/92

% Moisture: 3 decanted: (Y/N) N

Date Extracted: 07/07/92

Concentrated Extract Volume: 500.0 (uL)

Date Analyzed: 07/09/92

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 6.9

CONCENTRATION UNITS:

Number TICs found: 21

(ug/L or ug/Kg) UG/KG

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1.	UNKNOWN	6.40	210	J
2.	BLANK CONTAMINANT	6.55	210	BJ H ² O
3.	UNKNOWN	7.10	270	J'
4.	UNKNOWN HYDROCARBON	8.47	790	J
5.	UNKNOWN HYDROCARBON	9.44	650	J
6.	UNKNOWN HYDROCARBON	10.35	760	J
7.	UNKNOWN HYDROCARBON	11.19	620	J
8.	UNKNOWN HYDROCARBON	11.99	620	J
9.	UNKNOWN HYDROCARBON	13.47	690	J
10.	UNKNOWN HYDROCARBON	14.14	620	J
11.	UNKNOWN HYDROCARBON	14.79	450	J
12.	UNKNOWN HYDROCARBON	15.40	410	J
13.	UNKNOWN HYDROCARBON	15.99	380	J
14.	UNKNOWN HYDROCARBON	16.55	310	J
15.	UNKNOWN HYDROCARBON	17.10	380	J
16.	UNKNOWN HYDROCARBON	17.67	310	J
17.	UNKNOWN HYDROCARBON	18.30	310	J
18.	UNKNOWN HYDROCARBON	19.82	310	J
19.	UNKNOWN	22.15	340	J
20.	UNKNOWN HYDROCARBON	25.09	510	J
21.	UNKNOWN HYDROCARBON	19.00	140	J

1F
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

ORIGIN
(Ref)
CKY36

Lab Name: COMPUCHEM, RTP Contract: 68D10083

Lab Code: COMPU Case No.: 18347 SAS No.: _____ SDG No.: CKY28

Matrix: (soil/water) SOIL

Lab Sample ID: 499470

Sample wt/vol: 30.0 (g/mL) G

Lab File ID: GR099470A57

Level: (low/med) LOW

Date Received: 06/24/92

% Moisture: 10 decanted: (Y/N) N

Date Extracted: 07/07/92

Concentrated Extract Volume: 500.0 (uL)

Date Analyzed: 07/09/92

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 7.0

CONCENTRATION UNITS:

Number TICs found: 16

(ug/L or ug/Kg) UG/KG

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1.	UNKNOWN	6.42	150	J
2.	BLANK CONTAMINANT	6.48	150	BJ
3.	BLANK CONTAMINANT	6.57	220	BJ
4.	UNKNOWN	15.15	180	J
5.	UNKNOWN HYDROCARBON	17.10	110	J
6.	UNKNOWN	17.90	110	J
7.	UNKNOWN HYDROCARBON	18.30	110	J
8.	UNKNOWN HYDROCARBON	19.84	370	J
9.	UNKNOWN	19.95	330	J
10.	UNKNOWN HYDROCARBON	20.80	180	J
11.	UNKNOWN	21.34	1600	J
12.	UNKNOWN	22.12	1600	J
13.	UNKNOWN	23.37	150	J
14.	UNKNOWN	24.15	370	J
15.	UNKNOWN HYDROCARBON	25.12	1700	J
16.	UNKNOWN	25.57	110	J

1F
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

ORIGINAL
(Red) EPA SAMPLE NO.

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

CKY37

Lab Code: COMPU

Case No.: 18347

SAS No.: _____

SDG No.: CKY28

Matrix: (soil/water) SOIL

Lab Sample ID: 499471

Sample wt/vol: 30.0 (g/mL) G

Lab File ID: GH099471A15

Level: (low/med) LOW

Date Received: 06/24/92

% Moisture: 8 decanted: (Y/N) N

Date Extracted: 06/26/92

Concentrated Extract Volume: 500.0 (uL)

Date Analyzed: 07/02/92

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 6.1

CONCENTRATION UNITS:

Number TICs found: 17

(ug/L or ug/Kg) UG/KG

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1.	UNKNOWN	5.07	330	J
2.	BLANK-CONTAMINANT	5.17	1000	BJ
3.	BLANK-CONTAMINANT	5.22	140	BJ
4.	BLANK-CONTAMINANT	5.27	220	BJ
5.	BLANK-CONTAMINANT	5.32	360	BJ
6.	UNKNOWN	5.38	72	J
7.	UNKNOWN	5.60	220	J
8.	BLANK-CONTAMINANT	5.77	290	BJ
9.	UNKNOWN	5.98	140	J
10.	UNKNOWN	11.24	290	J
11.	UNKNOWN	12.85	110	J
12.	UNKNOWN	13.27	290	J
13.	LABORATORY-ARTIFACT	13.34	8300	J
14.	UNKNOWN	14.00	180	J
15.	UNKNOWN	17.65	330	J
16.	UNKNOWN	17.84	220	J
17.	UNKNOWN	27.42	510	J

1F
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

ORIGINAL
ReCKY38

Lab Name: COMPUCHEM, RTP Contract: 68D10083

Lab Code: COMPU Case No.: 18347 SAS No.: _____ SDG No.: CKY28

Matrix: (soil/water) SOIL

Lab Sample ID: 499473

Sample wt/vol: 30.0 (g/mL) G

Lab File ID: GR099473A57

Level: (low/med) LOW

Date Received: 06/24/92

% Moisture: 14 decanted: (Y/N) N

Date Extracted: 07/07/92

Concentrated Extract Volume: 500.0 (uL)

Date Analyzed: 07/09/92

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 6.4

CONCENTRATION UNITS:

Number TICs found: 7

(ug/L or ug/Kg) UG/KG

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1.	UNKNOWN	16.90	39	J
2.	UNKNOWN HYDROCARBON	18.30	120	J
3.	UNKNOWN HYDROCARBON	19.82	190	J
4.	UNKNOWN	21.29	430	J
5.	UNKNOWN HYDROCARBON	21.95	660	J
6.	UNKNOWN	24.14	350	J
7.	UNKNOWN HYDROCARBON	25.07	660	J

ORIGINAL
(Red)

Appendix E
Organic Regional Data Assessment Summary

TPO: ACTION FYIRegion IIIORGANIC REGIONAL DATA ASSESSMENT SUMMARY

CASE NO: 18347
 SDG NO: CKY14
 SOW: 3/90
 NO. OF SAMPLES: TWELVE (12)

LABORATORY: COMPU
 DATA USER: TOM BENNETT
 REVIEW COMPLETION DATE: 10/1/1992
 MATRIX: WATER

REVIEWER: ESAT

	VOA	BNA	PEST/PCB
1. HOLDING TIMES	<u>O</u>	<u>O</u>	<u>O</u>
2. GC-MS TUNE/GC PERFORMANCE	<u>O</u>	<u>O</u>	<u>O</u>
3. INITIAL CALIBRATIONS	<u>X</u>	<u>XA</u>	<u>O</u>
4. CONTINUING CALIBRATION	<u>X</u>	<u>X</u>	<u>O</u>
5. FIELD BLANKS (F=NOT APPLICABLE)	<u>X</u>	<u>X</u>	<u>O</u>
6. LABORATORY BLANKS	<u>X</u>	<u>X</u>	
7. SURROGATES	<u>O</u>	<u>X</u>	<u>X</u>
8. MATRIX SPIKE/DUPLICATES	<u>O</u>	<u>O</u>	
9. REGIONAL QC (F=NOT APPLICABLE)	<u>F</u>	<u>F</u>	<u>F</u>
10. INTERNAL STANDARDS	<u>O</u>	<u>O</u>	<u>O</u>
11. COMPOUND IDENTIFICATION	<u>O</u>	<u>O</u>	<u>O</u>
12. COMPOUND QUANTITATION	<u>O</u>	<u>O</u>	<u>O</u>
13. SYSTEM PERFORMANCE	<u>O</u>	<u>X</u>	<u>O</u>
14. OVERALL ASSESSMENT	<u>X</u>	<u>XA</u>	<u>X</u>

O = No problems or minor problems that do not affect data usability
 X = No more than about 5% of the data points are qualified as either estimated or unusable.

M = More than about 5% of the data points are qualified as estimated.

Z = More than about 5% of the data points are qualified as unusable.

A = TPO action requested; use in conjunction with one of the above codes.

TPO ACTION ITEMS: (See Item 3B.)AREAS OF CONCERN: (Documentation Attached, See Following Pages.)

TPO: ACTION FYI

Region III

ORGANIC REGIONAL DATA ASSESSMENT SUMMARY

CASE NO: 18347
 SDG NO: CKY28
 SOW: 3/90
 NO. OF SAMPLES: ELEVEN (11)

LABORATORY: COMPU
 DATA USER: TOM BENNETT
 REVIEW COMPLETION DATE: 10/1/1992
 MATRIX: SOIL

REVIEWER: ESAT

	VOA	BNA	PEST/PCB
1. HOLDING TIMES	<u>O</u>	<u>M</u>	<u>O</u>
2. GC-MS TUNE/GC PERFORMANCE	<u>O</u>	<u>O</u>	<u>O</u>
3. INITIAL CALIBRATIONS	<u>X</u>	<u>X</u>	<u>O</u>
4. CONTINUING CALIBRATION	<u>X</u>	<u>X</u>	<u>O</u>
5. FIELD BLANKS (F=NOT APPLICABLE)	<u>O</u>	<u>X</u>	<u>X</u>
6. LABORATORY BLANKS	<u>X</u>	<u>X</u>	<u>O</u>
7. SURROGATES	<u>O</u>	<u>XA</u>	<u>X</u>
8. MATRIX SPIKE/DUPLICATES	<u>O</u>	<u>O</u>	<u>O</u>
9. REGIONAL QC (F=NOT APPLICABLE)	<u>F</u>	<u>F</u>	<u>F</u>
10. INTERNAL STANDARDS	<u>O</u>	<u>O</u>	<u>O</u>
11. COMPOUND IDENTIFICATION	<u>O</u>	<u>O</u>	<u>O</u>
12. COMPOUND QUANTITATION	<u>O</u>	<u>O</u>	<u>O</u>
13. SYSTEM PERFORMANCE	<u>O</u>	<u>O</u>	<u>O</u>
14. OVERALL ASSESSMENT	<u>X</u>	<u>MA</u>	<u>X</u>

O = No problems or minor problems that do not affect data usability
 X = No more than about 5% of the data points are qualified as either estimated or unusable.

M = More than about 5% of the data points are qualified as estimated.

Z = More than about 5% of the data points are qualified as unusable.

A = TPO action requested; use in conjunction with one of the above codes.

TPO ACTION ITEMS: (See Item 7B.)AREAS OF CONCERN: (Documentation Attached, See Following Pages.)

(ed)
ORIGINAL

ORGANIC REGIONAL DATA ASSESSMENT SUMMARY NOTES
CASE: 18347 SDG: CKY14 AQUEOUS SAMPLES

Item 3A and 3B Several compounds failed precision criteria in the volatile and semivolatile initial calibrations. One target compound (3-nitroaniline) in the semivolatile fraction had a relative response factor (RRF) less than 0.05 in the initial calibration dated 6/15/92. (See Table I in Appendix F).

Item 4A and 4B Several compounds failed the precision criteria in the volatile semivolatile continuing calibrations. (See Table I in Appendix F.)

**Item 5A
5B
and 5C** The maximum concentrations of all compounds found in the analyses of field blanks are listed below:

<u>COMPOUND</u>	<u>CONCENTRATION</u>
methylene chloride*	22 µg/L
acetone*	15 µg/L
butylbenzylphthalate*	120 J µg/Kg
bis(2-ethylhexyl)phthalate*	2 J µg/L
methoxychlor	0.59 J µg/Kg

* Common Laboratory Contaminant

Item 6A and 6B The maximum concentrations of all compounds found in the analyses of laboratory method blanks are listed below:

<u>COMPOUND</u>	<u>CONCENTRATION</u>
methylene chloride*	30 µg/L
acetone*	31 µg/L
phenol	3 J µg/Kg
diethylphthalate*	1 J µg/Kg

* Common Laboratory Contaminant

Item 7B and 7C In the semivolatile analyses, aqueous samples CKY16, CKY19 and CKY20 had one (1) and sample CKY18 had two (2) (one acid and one base) surrogate recoveries above the QC limits. Sample CKY18MSD did not show any surrogate recovery due to an extraction error. In the pesticide/PCB analyses, aqueous samples CKY14 and CKY26 had one (1) each of their surrogate recoveries below the QC limits but greater than 10%. (See FORM II SV-1, case narrative and FORM II PEST-1 in Appendix F.)

Item 8B During the semivolatile analysis, sample CKY18MSD failed the spike recovery criteria due to an extraction error and no reanalysis was performed. (See FORM III SV-1 in Appendix F.)

ORIGINAL
(Red)

ORGANIC REGIONAL DATA ASSESSMENT SUMMARY NOTES
CASE: 18347 SDG: CKY28 SOIL SAMPLES

Item 1B The semivolatile extractions of samples CKY28, CKY31, CKY34, CKY35, CKY36 and CKY38 were performed thirteen (13) and sample CKY29 was performed seventeen (17) days from the date of sample collection. Although no technical holding time for the semivolatile extraction of the soil samples has been established, the technical holding time of seven (7) days for the aqueous samples has been exceeded by six (6) days and ten (10) days in these samples. The aqueous sample holding time was applied to these samples. (See Traffic Report in Appendix F.)

Item 3A and 3B Several compounds failed the precision criteria in the volatile and semivolatile initial calibrations. (See Table I in Appendix F).

Item 4A and 4B Several compounds failed the precision criteria in the volatile and semivolatile continuing calibrations. (See Table I in Appendix F.)

Item 5A 5B and 5C The maximum concentrations of all compounds found in the analyses of field blanks are listed below:

<u>COMPOUND</u>	<u>CONCENTRATION</u>
methylene chloride*	22 µg/L
acetone*	15 µg/L
butylbenzylphthalate*	120 J µg/Kg
bis(2-ethylhexyl)phthalate*	2 J µg/L
methoxychlor	0.59 J µg/Kg

* Common Laboratory Contaminant

Item 6A and 6B The maximum concentrations of all compounds found in the analyses of laboratory method blanks are listed below:

<u>COMPOUND</u>	<u>CONCENTRATION</u>
methylene chloride*	30 µg/L
acetone*	31 µg/L
phenol	3 J µg/Kg
diethylphthalate*	1 J µg/Kg

* Common Laboratory Contaminant

Item 7B and 7C The semivolatile analysis of sample CKY32 had the recovery for the acid surrogate, 2,4,6-tribromophenol less than 10%. Samples CKY32MS, CKY32MSD and CKY34 had one (1) of their surrogate recoveries outside the QC limits. In the pesticide/PCB analyses, samples CKY30, CKY32 and CKY36 had their surrogate recoveries below the QC limits but greater than 10% for

ORIGINAL
(Red)

ORGANIC REGIONAL DATA ASSESSMENT SUMMARY NOTES
CASE: 18347 SDG: CKY28 SOIL SAMPLES

tetrachloro-m-xylene (TCX) on one column. (See FORM II SV-2 and FORM II PEST-1 in Appendix F.)

- Item 8B The matrix spike duplicate recovery for sample CKY32 was above the QC limit for 4-nitrophenol in the semivolatile analyses. (See Form III SV-2 in Appendix F.)

ORIGINAL
(Red)

Appendix F
Support Documentation

TABLE I

Page | 6

ENVIRONMENTAL PROTECTION AGENCY REGION III
CALIBRATION CUTLERS
VOLATILE HSL COMPOUNDS
CONTRACTOR E.S.P.T.

ORIGINAL
(Red)

CASE/SER. NO. 18347

* See last page of this table for DEFINITION OF CODES.

TABLE I

page 2 of 14

ORIGINAL
(Red)

ENVIRONMENTAL PROTECTION AGENCY REGION III
CALIBRATION OUTLIERS
VOLATILE HSL COMPOUNDS
CONTRACTOR E.S.T.

CASE/SAS No. 18242

Instrument#	Init.	Cal.	Cont.	Cal.	Cont.	Cal.	Cont.	Cal.	Cont.	Cal.
DATE/TIME:	11-11-82 10:00 AM	RF 180								
Chloromethane										
Bromomethane										
Vinyl Chloride										
Chloroethane										
Methylene Chloride										
Acetone										
Carbon Disulfide										
1,1-Dichloroethene										
1,1-Dichloroethane										
Total 1,2-Dichloroethene										
Chloroform										
1,2-Dichloroethane										
2-Butanone										
1,1,1-Trichloroethane										
Carbon Tetrachloride										
Vinyl Acetate										
Bromodichloromethane										
1,2-Dichloropropane										
cis-1,3-Dichloropropene										
Dichloroethene										
Dibromochloromethane										
1,1,2-Trichloroethane										
Benzene										
trans-1,3-Dichloropropene										
Bromoform										
4-Methyl-2-Pentanone										
2-Hexanone										
Tetrachloroethene										
1,1,2-Tetrachloroethane										
Toluene										
Chlorobenzene										
Ethylbenzene										
Styrene										
Total Xylenes										
AFFECTED SAMPLES:										
Reviewer										
Initials/Date:	<u>H/a/2b2</u>									

* See last page of this table for DEFINITION OF CODES.

ENVIRONMENTAL PROTECTION AGENCY REGION III
CALIBRATION OUTLIERS
VOLATILE HSL COMPOUNDS
CONTRACTOR E&P

ORIGINAL
(Red)

CASE/SAS No. 18347

Instrument#	Init.	Cal.	Cont.	Cal.	Cont.	Cal.	Cont.	Cal.	Cont.	Cal.
DATE/TIME:										
	RF 18RSD	RF 18G	RF 18D							
Chloromethane										
Bromomethane										
Vinyl Chloride										
Chloroethane										
Methylene Chloride										
Acetone										
Carbon Disulfide										
1,1-Dichloroethene										
1,1-Dichloroethane										
Total-1,2-Dichloroethene										
Chloroform										
1,2-Dichloroethane										
2-Butanone										
1,1,1-Trichloroethane										
Carbon Tetrachloride										
Vinyl Acetate										
Bromodichloromethane										
1,2-Dichloropropane										
cis-1,3-Dichloropropene										
Trichloroethene										
Dibromochloromethane										
1,1,2-Trichloroethane										
Benzene										
trans-1,3-Dichloropropene										
Bromoform										
4-Methyl-2-Pentanone										
2-Hexanone										
Tetrachloroethene										
1,1,2-Tetrachloroethane										
luene										
Chlorobenzene										
Ethylbenzene										
Styrene										
Total Xylenes										
	CKV14	CKV18								
	CKV17	CKV26								
	CKV18MS									
	CKV18MSD									
AFFECTED	CKV19									
SAMPLES:	CKV20									
	CKV22									
Reviewer	CKV23									
Initials/Date:	CKV24									
	CKV25									

* See last page of this table for DEFINITION OF CODES.

TABLE I

page 4 of 14 ORIGINAL
(Red)

ENVIRONMENTAL PROTECTION AGENCY REGION III
CALIBRATION OUTLINES
SEMIVOLATILE HSL COMPOUNDS (Part 1 of 2)
CONTRACTOR ESAT

CASE / S.S. NO. 18347

* See last page of this table for DEFINITION OF CODES.

TABLE I

ORIGINAL
page 5 of 14 (Rev)

ENVIRONMENTAL PROTECTION AGENCY REGION III
CALIBRATION OUTLIERS
SEMI-VOLATILE HSL COMPOUNDS (Part 2 of 2)

CASE/ENR No 18347CONTRACTOR ESAT

Instrument#	Init.	Cal.	Cone.	Cal.	Cone.	Cal.	Cone.	Cal.	Cone.	Cal.
DATE/TIME:	11/25/92	03/04/92	07/15/92	07/15/92						
	IRF 182SD	*	IRF 18D	*	IRF 18D	*	IRF 18D	*	IRF 18D	*
Dibenzofuran										
2,4-Dinitrotoluene										
Diethylphthalate							-28.7°C			
4-Chlorophenyl-chenvlether										
Fluorene										
4-Nitroaniline										
1,6-Dinitro-2-methylbenzene										
N-Nitrosodiphenylamine										
4-Bromophenyl-chenvlether										
Hexachlorobenzene										
Pentachlorophenol										
Phenanthrene										
Anthracene										
Di-n-butylphthalate							-24.5°C			
Fluoranthene										
Pyrene										
Eucarbonylbenzylchloralate										
3,3-Dichlorocyanidine						-35.5°C	-60.0°C			
Benzo(a)anthracene										
Chrysene										
bis(2-Ethylhexyl)chloralate										
Di-n-octylphthalate										
Benzo(b)fluoranthene										
Benzo(k)fluoranthene										
Benzo(a)pyrene										
Indeno[1,2,3-d]pyrene										
dibenz(a,h)anthracene										
benzo(a,b,t)perylene										
<hr/>										
AFFECTED SAMPLES:	all sample C KY34 C KY28 on this page C KY31 C KY35 C KY36 C KY38									
Reviewer Initials/Date: <u>HP 9/10/92</u>										
<hr/>										

* See last page of this table for DEFINITION OF CODES.

TABLE I

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ENVIRONMENTAL PROTECTION AGENCY REGION III
CALIBRATION OUTLIERS
SEMICVOLATILE HSL COMPOUNDS (Part 1 of 2)

CASE/SAS No. 18347CONTRACTOR ESAT

Instrument	Init.	Cal.	Cont.	Cal.	Cont.	Cal.	Cont.	Cal.	Cont.	Cal.
DATE/TIME:	12/14/92	12/18								
Phenol										
bis(2-Chloroethyl)ether										
2-Chlorophenol										
1,3-Dichlorobenzene										
1,4-Dichlorobenzene										
Benzyl alcohol										
1,2-Dichlorobenzene										
4-Methylphenol										
bis(2-Chloroisopropyl)ether					-80.2C					
4-Methylphenol										
N-Nitroso-di-n-propylamine										
Hexachlorobutane										
Nitrobenzene										
Isophorone										
4-Nitrophenol										
4,4'-Dimethyldiphenol										
Benzoic acid										
bis(2-Chloroethyl)methane										
2,4-Dichlorophenol										
1,3,4-Trichlorobenzene										
Naphthalene										
4-Chloroaniline		37.1 C								
Hexachlorobutadiene										
4-Chloro-2-Methylphenol										
2-Methylnaphthalene										
Hexachlorocyclohexadiene										
4,6-Dichlorophenol										
2,4,5-Trichlorophenol										
2-Chloronaphthalene										
2-Nitroaniline				-28.2 C						
Dimethylphthalate										
Acenaphthylene										
2,6-Dinitrotoluene				-37.8 C						
3-Nitroaniline										
Acenaphthene										
2,4-Dinitrophenol				-43.2 C						
4-Nitrophenol				157.2 C						
AFFECTED SAMPLES:										
	all sample	CK430								
	on this page	CK432								
		CK432MS								
		CK432MD								
		CK433								
		CK437								

* See last page of this table for DEFINITION OF CODES.

TABLE I

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ORIGINAL
(Red)

ENVIRONMENTAL PROTECTION AGENCY REGION III
CALIBRATION CUTLERS
SEMITRIVOLATILE ESL COMPOUNDS (Part 2 of 2)

CASE/ENO No. 18342CONTRACTOR ESAT

Instrument#	Init.	Cal.	Conc.	Cal.	Conc.	Cal.	Conc.	Cal.	Conc.	Cal.
CWATIS	14.5/11/92	100%								
DATE/TIME:				IRF ISD	*	IRF ISD	*	IRF ISD	*	IRF ISD
Biphenolfuran										
2,4-Dinitrotoluene										
Diethylphthalate										
4-Chlorophenyl-chenvlether										
Fluorene										
4-Nitroaniline										
4,6-Dinitro-2-methyphenol										
Nitrosodiphenylamine										
4-Bromophenyl-chenvlether										
Hexachlorobenzene										
Pentachlorophenol										
Phenanthrene										
Anthracene										
Di-n-octylphthalate										
Fluoranthene										
Pyrene										
Ethylenephthalate										
3,3'-Dichlorobenzidine										
Benzo(a)anthracene										
Chrysene										
bis(2-Ethylhexyl)phthalate										
Di-n-octylphthalate										
Benzo(b)fluoranthene										
Benzo(k)fluoranthene										
Benzo(a)pyrene										
Indeno[1,2,3-cd]pyrene										
Benzo[a,h]anthracene										
Benzo[c,h,i]perylene										
Carboxylic										
AFFECTED SAMPLES:										
	all sample	CKY30								
	on this page	CKY32								
		CKY32MS								
		CKY32MS								
		CKY33								
		CKY37								

* See last page of this table for DEFINITION OF CODES.

TABLE I

page 8 of 14

ORIGINAL
(Red)

ENVIRONMENTAL PROTECTION AGENCY REGION III
CALIBRATION OUTLIERS
SEMIVOLATILE HSL COMPOUNDS (Part 1 of 2)

CASE/SAC No. 18347CONTRACTOR ESAT

Instrument#	OWA05	Int.	Cal.	Cont.	Cal.	Cont.	Cal.	Cont.	Cal.	Cont.	Cal.
DATE/TIME:		06/17/92	07/17/92								
		RF 14RSD	*	RF 14D	*	RF 14C	*	RF 14D	*	RF 14D	*
Phenol											
bis(2-Chloroethyl)ether											
2-Chlorophenol											
1,3-Dichlorobenzene											
1,4-Dichlorobenzene											
Benzyl alcohol											
2-Dichlorobenzene											
4-Methylphenol											
bis(2-Chloroisopropyl)ether											
4-Methylphenol											
N-Nitroso-di-n-propylamine											
Hexachloroethane											
Nitrobenzene											
Isochorone											
2-Nitrophenol											
2,4-Dimethoxybenzene											
Benzoic acid											
bis(2-Chloroethyl)ether											
2,4-Dichlorophenol											
1,2,4-Trichlorobenzene											
Naphthalene											
4-Chloraniline											
Hexachlorobutadiene											
4-Chloro-3-Methylphenol											
2-Methylnaphthalene											
Hexachlorobutadiene											
4,6-Trichlorophenol											
2,4,6-Trichlorophenol											
2-Chloronaphthalene											
2-Nitroaniline											
Dimethylchloralate											
Acenaphthylene											
2,6-Dinitrotoluene											
3-Nitroaniline											
Acenaphthene											
2,4-Dinitrophenol											
4-Nitrophenol											
AFFECTED SAMPLES:											
Reviewer											
Initials/Date:	HP	9/10/92									

all sample SK429
on this page

* See last page of this table for DEFINITION OF CODES.

TABLE I

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ORIGINAL
Reed

ENVIRONMENTAL PROTECTION AGENCY REGION III
CALIBRATION OUTLIERS
SEMOVOLATILE HSL COMPOUNDS (Part 1 of 2)

CASE/SAS No. 18347CONTRACTOR ESAT

Chemical	Init.	Cal.	Cont.	Cal.	Cont.	Cal.	Cont.	Cal.	Cont.	Cal.
CASE DATE:	10/26/85	10/26/85	10/26/85	10/26/85	10/26/85	10/26/85	10/26/85	10/26/85	10/26/85	10/26/85
Phenol										
bis(2-Chloroethyl)ether										
2-Chlorophenol										
1,3-Dichlorobenzene										
1,4-Dichlorobenzene										
Benzyl alcohol										
1,6-Dichlorobenzene										
4-Etoxyphenol										
bis(2-Chloroisopropenyl)ether										
3-Methoxyphenol										
N-Nitroso-di-n-propylamide										
Hexachloroethane										
Nitrobenzene										
Isocyanone										
2-Nitrophenol										
2,4-Dinitrophenol										
Benzoic acid										
bis(2-Chloroethyl)benzene										
2,4-Dichlorophenol										
1,3,5-Trichlorobenzene										
Naphthalene										
3-Chloronaphthalene		33.9			-44.2	C				
Hexachlorocyclohexene										
2-Chloro-3-Methoxyphenol										
2-Methoxynaphthalene										
Acidchlorocyclopentadiene										
2,4-Dichlorophenol										
2,4,5-Trichlorophenol										
3-Chlorodibenzofuran										
3-Nitroaniline		6.044	89.7	F	-3523	C				
Acenaphthene										
2,6-Dinitrotoluene										
3-Nitroaniline										
Acenaphthene										
2,4-Dinitrophenol										
4-Nitrophenol										
AFFECTED SAMPLES:										
Reviewer										
Initials/Date:	<u>HP 9/9/92</u>									

all samples C K 724
this page C K 725
C K 726

* See last page of this table for DEFINITION OF CODES.

TABLE I

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(Red) ORIGINAL

ENVIRONMENTAL PROTECTION AGENCY REGION III
CALIBRATION OUTLINES
SEMIVOLATILE HSL COMPOUNDS (Part 2 of 2)
CONTRACTOR E-SAT

CASE/SAC No. 18347

* See last page of this table for DEFINITION OF CODES.

TABLE I

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ENVIRONMENTAL PROTECTION AGENCY REGION III
 CALIBRATION OUTLINES
 SEMIVOLATILE HSL COMPOUNDS (Part 2 of 2)

ORIGINAL
(Red)CASE/SMS No. 18347CONTRACTOR ESAT

Instrument#	Init.	Cal.	Cont.	Cal.	Cont.	Cal.	Cont.	Cal.	Cont.	Cal.
DATE/TIME:	10/21/92 7:45	10/21/92 7:45	10/21/92 7:45	10/21/92 7:45	10/21/92 7:45	10/21/92 7:45	10/21/92 7:45	10/21/92 7:45	10/21/92 7:45	10/21/92 7:45
Dibenzofuran										
2,4-Dinitrotoluene										
Diethylphthalate										
4-Chlorophenyl-phenylether										
Fluorene										
4-Nitroaniline		64.3 I		-46.1 C						
3,6-Dinitro-2-methylphenol										
N-Nitrosodiphenylamine										
4-Bromophenyl-phenylether										
Hexachlorobenzene										
Pentachlorophenol										
Phenanthrene										
Anthracene										
Di-n-butylphthalate										
Fluoranthene										
Biphenyl										
Butylbenzylphthalate										
3,3'-Dichlorobenzidine		70.9 I		-95.2 C						
benzo(a)anthracene										
Chrysene										
bis(2-Ethylhexyl)phthalate										
Di-n-octylphthalate										
benzo(b)fluoranthene										
benzo(k)fluoranthene										
benzo(a)pyrene										
Indeno[1,2,3-cd]pyrene										
Dibenz(a,h)anthracene										
benzo(a,h,i)perylene										
Carbazole										
AFFECTED SAMPLES:										
	all samples on CKY24									
	this page	CKY25								
		CKY26								

Reviewer
 Initials/Date: HP 9/9/92

* See last page of this table for DEFINITION OF CODES.

TABLE I

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ORIGINAL
(Red)

ENVIRONMENTAL PROTECTION AGENCY REGION III
CALIBRATION CUTTERS
SEMICVOLATILE HSL COMPOUNDS (Part 1 of 2)

CASE/SAS No. 18347

CONTRACTOR ESCAT

Instrument	Cal.	Cont.	Cal.	Cont.	Cal.	Cont.	Cal.	Cont.	Cal.
DATE/TIME:	10/23/92	10/23/92	10/23/92	10/23/92	10/23/92	10/23/92	10/23/92	10/23/92	10/23/92
Phenol	-	-	-	-	-	-	-	-	-
bis(2-Chloroethyl)ether	-	-	-	-	-	-	-	-	-
2-Chlorostyrene	-	-	-	-	-	-	-	-	-
1,3-Dichlorobenzene	-	-	-	-	-	-	-	-	-
1,4-Dichlorobenzene	-	-	-	-	-	-	-	-	-
Benzyl alcohol	-	-	-	-	-	-	-	-	-
2-Chlorobenzene	-	-	-	-	-	-	-	-	-
o-Methylstyrene	-	-	-	-	-	-	-	-	-
bis(2-Chloroisopropyl)ether	-	-	-	-	-	-	-	-	-
o-Methylstyrene	-	-	-	-	-	-	-	-	-
N-Nitroso-di-n-propylamine	-	-	-	-	-	-	-	-	-
Hexachloroethane	-	-	-	-	-	-	-	-	-
Nicotinamide	-	-	-	-	-	-	-	-	-
Isophorone	-	-	-	-	-	-	-	-	-
2-Nitropropanoic acid	-	-	-	-	-	-	-	-	-
2,4-Dichlorobenzene	-	-	-	-	-	-	-	-	-
Naphthalene	-	-	-	-	-	-	-	-	-
o-Chlorostyrene	-	-	-	-	-	-	-	-	-
Hexachlorocyclohexene	-	-	-	-	-	-	-	-	-
2-Chloro-3-Methylbenzene	-	-	-	-	-	-	-	-	-
2-Chlorvinylbenzaldehyde	-	-	-	-	-	-	-	-	-
1,3-Chloro-5-vinylbenzene	-	-	-	-	-	-	-	-	-
2,4,6-Trichlorobenzene	-	-	-	-	-	-	-	-	-
2,4,5-Trichlorobenzene	-	-	-	-	-	-	-	-	-
2-Chloronaphthalene	-	-	-	-	-	-	-	-	-
3-Nitroaniline	-	-	-	-	-	-	-	-	-
Dimethylchloralate	-	-	-	-	-	-	-	-	-
Acenaphthylene	-	-	-	-	-	-	-	-	-
2,6-Dinitrotoluene	-	-	-	-	-	-	-	-	-
3-Nitroaniline	-	-	-	-	-	-	-	-	-
Acenaphthene	-	-	-	-	-	-	-	-	-
2,4-Dinitrophenol	-	-	-	-	-	-	-	-	-
4-Nitrophenol	-	-	-	-	-	-	-	-	-
AFFECTED SAMPLES:	All samples in CKY10								
Reviewer Initials/Date:	HP	9/9/92	CKY20						
			CKY10MSD						
			CKY22						
			CKY23						
			CKY15						
			CKY16						
			CKY17						
			CKY18						
			CKY19						

* See last page of this table for DEFINITION OF CODES.

TABLE I

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ORIGINAL
(Red)

ENVIRONMENTAL PROTECTION AGENCY REGION III
CALIBRATION OUTLINES
SEMICOLATILE HSL COMPOUNDS (Part 2 of 2)

CASE/SAC No. 18347 CONTRACTOR ESAT

Instrument#	Int.	Cal.	Cont.	Cal.	Cont.	Cal.	Cont.	Cal.	Cont.	Cal.
DATE/TIME:	10/25/83(92)	10/25/83(92)								
Dibenzofuran	✓	✓	✓	✓	✓	✓	✓	✓	✓	✓
3,4-Dinitrotoluene	✓	✓	✓	✓	✓	✓	✓	✓	✓	✓
Diethylphthalate	✓	✓	✓	✓	✓	✓	✓	✓	✓	✓
4-Chlorophenyl-phenylether	✓	✓	✓	✓	✓	✓	✓	✓	✓	✓
Fluorine	✓	✓	✓	✓	✓	✓	✓	✓	✓	✓
4-Nitroaniline	✓	✓	✓	✓	✓	✓	✓	✓	✓	✓
4,6-Dinitro-2-methoxyphenol	✓	✓	✓	✓	✓	✓	✓	✓	✓	✓
Nitrosodiphenylamine	✓	✓	✓	✓	✓	✓	✓	✓	✓	✓
4-Bromophenyl-phenylether	✓	✓	✓	✓	✓	✓	✓	✓	✓	✓
Hexachlorobenzene	✓	✓	✓	✓	✓	✓	✓	✓	✓	✓
Pentachlorophenol	✓	✓	✓	✓	✓	✓	✓	✓	✓	✓
Phenanthrene	✓	✓	✓	✓	✓	✓	✓	✓	✓	✓
Anthracene	✓	✓	✓	✓	✓	✓	✓	✓	✓	✓
Di-n-octylphthalate	✓	✓	✓	✓	✓	✓	✓	✓	✓	✓
Fluoranthene	✓	✓	✓	✓	✓	✓	✓	✓	✓	✓
Pyrene	✓	✓	✓	✓	✓	✓	✓	✓	✓	✓
Butylbenzylphthalate				259 C						
3,3'-Dichlorobenzidine				252 C						
Benz(a)anthracene	✓	✓	✓	✓	✓	✓	✓	✓	✓	✓
Chrysene	✓	✓	✓	✓	✓	✓	✓	✓	✓	✓
Bis(2-Ethylhexyl)phthalate	✓	✓	✓	✓	✓	✓	✓	✓	✓	✓
Di-n-octylphthalate	✓	✓	✓	✓	✓	✓	✓	✓	✓	✓
Benz(b)fluoranthene	✓	✓	✓	✓	✓	✓	✓	✓	✓	✓
Benz(k)fluoranthene	✓	✓	✓	✓	✓	✓	✓	✓	✓	✓
Benz(a)pyrene	✓	✓	✓	✓	✓	✓	✓	✓	✓	✓
Diene(1,3,5-odd)octane	✓	✓	✓	✓	✓	✓	✓	✓	✓	✓
Benz(a)anthracene	✓	✓	✓	✓	✓	✓	✓	✓	✓	✓
Benz(c,h,i)perylene	✓	✓	✓	✓	✓	✓	✓	✓	✓	✓
Calibration:	all samples	CKY16								
AFFECTED SAMPLES:	1st page	CKY20								
		CKY14M								
Reviewer Initials/Date:	4P9/9/92	CKY14D								
		CKY22								
		CKY23								
		CKY15								
		CKY16								
		CKY17								
		CKY18								
		CKY19								

* See last page of this table for DEFINITION OF CODES.

DEFINITION OF CODES USED IN TABLE I

I = %RSD exceeded 30% in the initial calibration, positive results are qualified "J". When the %RSD exceeded 50%, quantitation limits are qualified "UJ".

C = %D exceeded 25% in the continuing calibration, positive results are qualified "J". When the %D exceeded 50%, quantitation limits are qualified "UJ".

F = RF less than 0.05 in the calibration. All quantitation limits are qualified "R".

+ = The "B" qualifier, denoting blank contamination, supersedes the qualifier issued in this table.

R = The "R" qualifier, denoting unusable results, supersedes the qualifier issued in this table.

6A
VOLATILE ORGANICS INITIAL CALIBRATION DATA

Lab Name: COMPUCHEM, RTP Contract: 68D10083 ORIGINAL
(Red)
 Lab Code: COMPU Case No.: 18347 SAS No.: SDG No.: CKY28
 Instrument ID: F50054 Calibration Date(s): 06/09/92 06/10/92
 Heated Purge: (Y/N): Y Calibration Times: 1803 0140
 GC Column: DB-624 ID: 0.530(mm)

LAB FILE ID:	RRF10 = <u>GT920610C54</u>	RRF20 = <u>GS920610C54</u>					
RRF50=	<u>GU920609B54</u>	<u>RRF100= GT920609B54</u>					
RRF200=	<u>GZ920609B54</u>						
COMPOUND	RRF10	RRF20	RRF50	RRF100	RRF200	RRF	% RSD
Chloromethane	1.381	1.037	0.997	0.891	0.986	1.058	17.8
Bromomethane	* 1.751	1.350	1.091	1.041	1.173	1.281	22.5*
Vinyl Chloride	* 1.572	1.361	1.172	1.117	1.150	1.274	15.0*
Chloroethane	0.967	0.869	0.714	0.655	0.719	0.785	16.4
Methylene Chloride	4.436	3.579	1.923	1.894	1.837	2.734	44.0
Acetone	0.843	0.874	0.577	0.498	0.406	0.640	32.7
Carbon Disulfide	4.978	4.699	3.983	3.999	3.822	4.296	11.9
1,1-Dichloroethene	* 1.645	1.657	1.439	1.509	1.395	1.529	7.8*
1,1-Dichloroethane	* 3.279	3.002	2.678	2.718	2.557	2.847	10.2*
1,2-Dichloroethene (total)	1.756	1.738	1.474	1.555	1.444	1.593	9.2
Chloroform	* 4.211	3.778	3.394	3.527	3.373	3.657	9.6*
1,2-Dichloroethane	* 2.411	2.294	1.953	2.017	1.969	2.129	9.8*
2-Butanone	0.429	0.491	0.464	0.521	0.425	0.466	8.8
1,1,1-Trichloroethane	* 0.750	0.692	0.658	0.652	0.659	0.682	6.0*
Carbon Tetrachloride	* 0.749	0.717	0.664	0.669	0.692	0.698	5.1*
Bromodichloromethane	* 0.828	0.801	0.750	0.758	0.776	0.783	4.1*
1,2-Dichloropropane	0.396	0.385	0.349	0.342	0.344	0.363	7.0
cis-1,3-Dichloropropene	* 0.533	0.551	0.528	0.542	0.558	0.542	2.3*
Trichloroethene	* 0.525	0.500	0.446	0.439	0.437	0.469	8.6*
Dibromochloromethane	* 0.697	0.668	0.650	0.658	0.660	0.667	2.7*
1,1,2-Trichloroethane	* 0.388	0.375	0.355	0.350	0.342	0.362	5.2*
Benzene	* 1.039	0.962	0.885	0.847	0.838	0.914	9.3*
Trans-1,3-Dichloropropene	* 0.316	0.344	0.355	0.380	0.398	0.359	8.9*
Bromoform	* 0.549	0.545	0.545	0.554	0.543	0.547	0.8*
4-Methyl-2-Pentanone	0.386	0.435	0.428	0.423	0.399	0.415	5.2
2-Hexanone	0.260	0.290	0.286	0.297	0.271	0.281	5.3
Tetrachloroethene	* 0.623	0.589	0.540	0.519	0.499	0.554	9.2*
1,1,2,2-Tetrachloroethane	* 0.818	0.798	0.774	0.751	0.715	0.771	5.2*
Toluene	* 1.523	1.388	1.269	1.222	1.206	1.322	10.1*
Chlorobenzene	* 1.187	1.123	1.024	0.994	0.978	1.061	8.5*
Ethylbenzene	* 0.527	0.492	0.460	0.453	0.445	0.475	7.1*
Styrene	* 1.033	0.958	0.911	0.902	0.892	0.939	6.2*
Xylene (total)	* 0.649	0.614	0.575	0.563	0.556	0.591	6.6*
Toluene-d8	1.271	1.137	1.107	1.061	1.049	1.125	7.9
Bromofluorobenzene	* 1.133	1.001	1.016	0.974	0.970	1.019	6.5*
1,2-Dichloroethane-d4	2.194	1.916	1.830	1.894	1.857	1.938	7.6

* Compounds with required minimum RRF and maximum %RSD values.
 All other compounds must meet a minimum RRF of 0.010.

7A
VOLATILE CONTINUING CALIBRATION CHECK

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU

Case No.: 18347

SAS No.: _____

SDG No.: CKY28

ORIGINAL

Instrument ID: F50054

Calibration date: 06/24/92 Time: 1346

Lab File ID: GS920624A54

Init. Calib. Date(s): 06/09/92 06/10/92

Heated Purge: (Y/N) Y

Init. Calib. Times: 1803 0140

GC Column: DB-624 ID: 0.530 (mm)

Samples

COMPOUND	RRF	RRF50	MIN RRF	%D	MAX %D	SAMPLE
Chloromethane	1.058	0.715	(32.4)			CKY28
Bromomethane	1.281	1.417	0.100	-10.6	25.0	CKY29
Vinyl Chloride	1.274	1.074	0.100	15.7	25.0	CKY30
Chloroethane	0.785	0.861		-9.7		
Methylene Chloride	2.734	2.344		14.3		CKY31
Acetone	0.640	0.609		4.8		
Carbon Disulfide	4.296	3.786		11.9		CKY34
1,1-Dichloroethene	1.529	1.574	0.100	-2.9	25.0	
1,1-Dichloroethane	2.847	2.700	0.200	5.2	25.0	CKY35
1,2-Dichloroethene (total)	1.593	1.575		1.1		CKY37
Chloroform	3.657	3.826	0.200	-4.6	25.0	
1,2-Dichloroethane	2.129	2.355	0.100	-10.6	25.0	
2-Butanone	0.466	0.431		7.5		
1,1,1-Trichloroethane	0.682	0.696	0.100	-2.1	25.0	
Carbon Tetrachloride	0.698	0.762	0.100	-9.2	25.0	
Bromodichloromethane	0.783	0.805	0.200	-2.8	25.0	
1,2-Dichloropropane	0.363	0.343		5.5		
cis-1,3-Dichloropropene	0.542	0.481	0.200	11.2	25.0	
Trichloroethene	0.469	0.465	0.300	0.9	25.0	
Dibromochloromethane	0.667	0.719	0.100	-7.8	25.0	
1,1,2-Trichloroethane	0.362	0.381	0.100	-5.2	25.0	
Benzene	0.914	0.888	0.500	2.8	25.0	
Trans-1,3-Dichloropropene	0.359	0.314	0.100	12.5	25.0	
Bromoform	0.547	0.604	0.100	-10.4	25.0	
4-Methyl-2-Pentanone	0.415	0.333		19.8		
2-Hexanone	0.281	0.211		24.9		
Tetrachloroethene	0.554	0.571	0.200	-3.1	25.0	
1,1,2,2-Tetrachloroethane	0.771	0.755	0.500	2.1	25.0	
Toluene	1.322	1.272	0.400	3.8	25.0	
Chlorobenzene	1.061	1.006	0.500	5.2	25.0	
Ethylbenzene	0.475	0.468	0.100	1.5	25.0	
Styrene	0.939	0.882	0.300	6.1	25.0	
Xylene (total)	0.591	0.589	0.300	0.3	25.0	
Toluene-d8	1.125	1.098		2.4		
Bromofluorobenzene	1.019	1.002	0.200	1.7	25.0	
1,2-Dichloroethane-d4	1.938	2.184		-12.7		

All other compounds must meet a minimum RRF of 0.010.

7A
VOLATILE CONTINUING CALIBRATION CHECK

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

ORIGINAL
(Req)

Lab Code: COMPU Case No.: 18347 SAS No.: _____ SDG No.: CKY28

Instrument ID: F50054 Calibration date: 06/25/92 Time: 0507

Lab File ID: GS920625C54 Init. Calib. Date(s): 06/09/92 06/10/92

Heated Purge: (Y/N) Y Init. Calib. Times: 1803 0140

GC Column: DB-624 ID: 0.530 (mm)

COMPOUND	RRF	RRF50	MIN RRF	%D	MAX %D	Sample
Chloromethane	1.058	1.166		-10.2		CK432
Bromomethane	1.281	1.496	0.100	-16.8	25.0	CK432M
Vinyl Chloride	1.274	1.348	0.100	-5.8	25.0	CK433
Chloroethane	0.785	0.830		-5.7		CK436
Methylene Chloride	2.734	2.219		18.8		
Acetone	0.640	0.323		49.5		CK438
Carbon Disulfide	4.296	4.311		-0.3		
1,1-Dichloroethene	1.529	1.554	0.100	-1.6	25.0	
1,1-Dichloroethane	2.847	2.803	0.200	1.5	25.0	
1,2-Dichloroethene (total)	1.593	1.579		0.9		
Chloroform	3.657	3.828	0.200	-4.7	25.0	
1,2-Dichloroethane	2.129	2.267	0.100	-6.5	25.0	
2-Butanone	0.466	0.402		13.7		
1,1,1-Trichloroethane	0.682	0.700	0.100	-2.6	25.0	
Carbon Tetrachloride	0.698	0.780	0.100	-11.8	25.0	
Bromodichloromethane	0.733	0.822	0.200	-5.0	25.0	
1,2-Dichloropropane	0.363	0.347		4.4		
cis-1,3-Dichloropropene	0.542	0.482	0.200	11.1	25.0	
Trichloroethene	0.469	0.480	0.300	-2.3	25.0	
Dibromochloromethane	0.667	0.706	0.100	-5.8	25.0	
1,1,2-Trichloroethane	0.362	0.350	0.100	3.3	25.0	
Benzene	0.914	0.884	0.500	3.3	25.0	
Trans-1,3-Dichloropropene	0.359	0.315	0.100	12.3	25.0	
Bromoform	0.547	0.574	0.100	-4.9	25.0	
4-Methyl-2-Pentanone	0.415	0.323		22.2		
2-Hexanone	0.281	0.177		37.0		
Tetrachloroethene	0.554	0.572	0.200	-3.2	25.0	
1,1,2,2-Tetrachloroethane	0.771	0.655	0.500	15.0	25.0	
Toluene	1.322	1.267	0.400	4.2	25.0	
Chlorobenzene	1.061	1.022	0.500	3.7	25.0	
Ethylbenzene	0.475	0.461	0.100	2.9	25.0	
Styrene	0.939	0.906	0.300	3.5	25.0	
Xylene (total)	0.591	0.584	0.300	1.2	25.0	
Toluene-d8	1.125	1.104		1.9		
Bromofluorobenzene	1.019	1.024	0.200	-0.5	25.0	
1,2-Dichloroethane-d4	1.938	2.090		-7.8		

All other compounds must meet a minimum RRF of 0.010.

7A
VOLATILE CONTINUING CALIBRATION CHECK

ORIGINAL
(Red)

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 18347 SAS No.: _____ SDG No.: CKY28

Instrument ID: F50054 Calibration date: 06/26/92 Time: 0840

Lab File ID: GS920626A54 Init. Calib. Date(s): 06/09/92 06/10/92

Heated Purge: (Y/N) Y Init. Calib. Times: 1803 0140

GC Column: DB-624 ID: 0.530 (mm)

Samples

CKY32.M

COMPOUND	RRF	RRF50	MIN RRF	%D	%D
Chloromethane	1.058	0.989	6.5		
Bromomethane	1.281	1.417	0.100	-10.6	25.0
Vinyl Chloride	1.274	1.236	0.100	3.0	25.0
Chloroethane	0.785	0.892		-13.6	
Methylene Chloride	2.734	1.966		(28.1)	
Acetone	0.640	0.451		29.5	
Carbon Disulfide	4.296	3.836		10.7	
1,1-Dichloroethene	1.529	1.586	0.100	-3.7	25.0
1,1-Dichloroethane	2.847	2.743	0.200	3.7	25.0
1,2-Dichloroethene (total)	1.593	1.548		2.8	
Chloroform	3.657	3.830	0.200	-4.7	25.0
1,2-Dichloroethane	2.129	2.333	0.100	-9.6	25.0
2-Butanone	0.466	0.398		14.6	
1,1,1-Trichloroethane	0.682	0.719	0.100	-5.4	25.0
Carbon Tetrachloride	0.698	0.776	0.100	-11.2	25.0
Bromodichloromethane	0.783	0.832	0.200	-6.3	25.0
1,2-Dichloropropane	0.363	0.355		2.2	
cis-1,3-Dichloropropene	0.542	0.514	0.200	5.2	25.0
Trichloroethene	0.469	0.476	0.300	-1.5	25.0
Dibromochloromethane	0.667	0.730	0.100	-9.4	25.0
1,1,2-Trichloroethane	0.362	0.370	0.100	-2.2	25.0
Benzene	0.914	0.898	0.500	1.8	25.0
Trans-1,3-Dichloropropene	0.359	0.339	0.100	5.6	25.0
Bromoform	0.547	0.616	0.100	-12.6	25.0
4-Methyl-2-Pentanone	0.415	0.383		7.7	
2-Hexanone	0.281	0.238		15.3	
Tetrachloroethene	0.554	0.578	0.200	-4.3	25.0
1,1,2,2-Tetrachloroethane	0.771	0.763	0.500	1.0	25.0
Toluene	1.322	1.310	0.400	0.9	25.0
Chlorobenzene	1.061	1.059	0.500	0.2	25.0
Ethylbenzene	0.475	0.474	0.100	0.2	25.0
Styrene	0.939	0.932	0.300	0.7	25.0
Xylene (total)	0.591	0.618	0.300	-4.6	25.0
Toluene-d8	1.125	1.144		-1.7	
Bromofluorobenzene	1.019	1.041	0.200	-2.2	25.0
1,2-Dichloroethane-d4	1.938	2.130		-9.9	

All other compounds must meet a minimum RRF of 0.010.

6A
VOLATILE ORGANICS INITIAL CALIBRATION DATALab Name: COMPUCHEM, RTPContract: 68D10083Lab Code: COMPU Case No.: 18347

SAS No.: _____

SDG No.: CKY14Instrument ID: F50056Calibration Date(s): 06/22/9206/23/92Heated Purge: (Y/N): NCalibration Times: 15200056GC Column: DB624 ID: 0.530(mm)

LAB FILE ID:	RRF10 = <u>CU920622B56</u>	RRF20 = <u>CA920622A56</u>
RRF50=	<u>CT920622B56</u>	RRF100= <u>CY920622A56</u>
	RRF200= <u>CX920622A56</u>	

COMPOUND	RRF10	RRF20	RRF50	RRF100	RRF200	RRF	% RSD
Chloromethane	0.061	0.041	0.048	0.048	0.051	0.050	14.5
Bromomethane	* 1.063	1.079	1.006	0.964	0.913	1.005	6.9*
Vinyl Chloride	* 0.633	0.415	0.391	0.382	0.353	0.435	26.0*
Chloroethane	0.741	0.847	0.741	0.734	0.688	0.750	7.8
Methylene Chloride	2.875	3.490	1.514	1.477	1.292	2.130	46.4
Acetone	0.262	0.328	0.214	0.196	0.172	0.234	26.4
Carbon Disulfide	3.856	4.114	3.856	4.040	3.573	3.888	5.4
1,1-Dichloroethene	* 1.435	1.350	1.319	1.326	1.191	1.324	6.6*
1,1-Dichloroethane	* 2.833	3.054	2.839	2.981	2.791	2.910	3.7*
1,2-Dichloroethene (total)	1.470	1.596	1.480	1.523	1.397	1.494	4.9
Chloroform	* 2.733	2.980	2.811	2.847	2.709	2.816	3.8*
1,2-Dichloroethane	* 1.500	1.720	1.673	1.684	1.653	1.646	5.2*
2-Butanone	0.177	0.442	0.371	0.385	0.386	0.352	28.9
1,1,1-Trichloroethane	* 0.474	0.486	0.477	0.427	0.433	0.459	5.9*
Carbon Tetrachloride	* 0.441	0.439	0.452	0.411	0.417	0.432	4.0*
Bromodichloromethane	* 0.473	0.606	0.535	0.485	0.576	0.535	10.8*
1,2-Dichloropropane	0.377	0.461	0.420	0.388	0.430	0.415	8.1
cis-1,3-Dichloropropene	* 0.510	0.533	0.586	0.539	0.558	0.555	5.7*
Trichloroethene	* 0.480	0.457	0.453	0.423	0.410	0.445	6.3*
Dibromochloromethane	* 0.439	0.444	0.495	0.439	0.459	0.455	5.2*
1,1,2-Trichloroethane	* 0.333	0.342	0.359	0.299	0.313	0.329	7.2*
Benzene	* 1.022	1.156	1.019	0.926	0.929	1.010	9.3*
Trans-1,3-Dichloropropene	* 0.326	0.375	0.383	0.360	0.374	0.364	6.2*
Bromoform	* 0.322	0.319	0.378	0.327	0.350	0.339	7.3*
4-Methyl-2-Pentanone	0.149	0.204	0.192	0.197	0.196	0.188	11.7
2-Hexanone	0.149	0.204	0.192	0.197	0.196	0.188	11.7
Tetrachloroethene	* 0.539	0.515	0.508	0.484	0.437	0.497	7.8*
1,1,2,2-Tetrachloroethane	* 0.504	0.494	0.522	0.476	0.468	0.493	4.4*
Toluene	* 1.321	1.420	1.344	1.325	1.187	1.319	6.4*
Chlorobenzene	* 1.018	1.072	1.057	1.016	0.932	1.019	5.3*
Ethylbenzene	* 0.520	0.538	0.543	0.530	0.484	0.523	4.5*
Styrene	* 0.952	1.018	1.027	0.991	0.917	0.981	4.7*
Xylene (total)	* 0.629	0.662	0.657	0.638	0.581	0.633	5.1*
Toluene-d8	-	1.171	1.216	1.183	1.159	1.116	3.1
Bromofluorobenzene	* 0.789	0.827	0.809	0.803	0.813	0.808	1.7*
1,2-Dichloroethane-d4	1.431	1.632	1.608	1.629	1.718	1.604	6.6

* Compounds with required minimum RRF and maximum %RSD values.
All other compounds must meet a minimum RRF of 0.010.

FORM VI VOA

ORIGINAL
(Red)

3/90

162

SAMPLE DATA PACKAGE

10217 CKY 14

7A
VOLATILE CONTINUING CALIBRATION CHECK

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

ORIGINAL
(Red)

Lab Code: COMPU

Case No.: 18347

SAS No.: _____

SDG No.: CKY14

Instrument ID: F50056

Calibration date: 06/24/92 Time: 0648

Lab File ID: CS920624C56

Init. Calib. Date(s): 06/22/92

06/23/92

Heated Purge: (Y/N) N

Init. Calib. Times: 1520

0056

GC Column: DB624 ID: 0.530(mm)

COMPOUND	RRF	RRF50	MIN RRF	%D	MAX %D	Sample
Chloromethane	0.050	0.076		-52.0		CKY16
Bromomethane	1.005	0.936	0.100	6.9	25.0	
Vinyl Chloride	0.435	0.411	0.100	5.5	25.0	
Chloroethane	0.750	0.645		14.0		
Methylene Chloride	2.130	1.531		28.1		
Acetone	0.234	0.229		2.1		
Carbon Disulfide	3.883	3.596		7.5		
1,1-Dichloroethene	1.324	1.282	0.100	3.2	25.0	
1,1-Dichloroethane	2.910	2.855	0.200	1.9	25.0	
1,2-Dichloroethene (total)	1.494	1.516		-1.5		
Chloroform	2.816	2.884	0.200	-2.4	25.0	
1,2-Dichloroethane	1.646	1.723	0.100	-4.7	25.0	
2-Butanone	0.352	0.354		-0.6		
1,1,1-Trichloroethane	0.459	0.452	0.100	1.5	25.0	
Carbon Tetrachloride	0.432	0.406	0.100	6.0	25.0	
Bromodichloromethane	0.535	0.498	0.200	6.9	25.0	
1,2-Dichloropropane	0.415	0.406		2.2		
cis-1,3-Dichloropropene	0.555	0.566	0.200	-2.0	25.0	
Trichloroethene	0.445	0.469	0.300	-5.4	25.0	
Dibromochloromethane	0.455	0.466	0.100	-2.4	25.0	
1,1,2-Trichloroethane	0.329	0.356	0.100	-8.2	25.0	
Benzene	1.010	0.983	0.500	2.7	25.0	
Trans-1,3-Dichloropropene	0.364	0.373	0.100	-2.5	25.0	
Bromoform	0.339	0.353	0.100	-4.1	25.0	
4-Methyl-2-Pentanone	0.188	0.329		75.0		
2-Hexanone	0.188	0.180		4.3		
Tetrachloroethene	0.497	0.535	0.200	-7.6	25.0	
1,1,2,2-Tetrachloroethane	0.493	0.574	0.500	-16.4	25.0	
Toluene	1.319	1.402	0.400	-6.3	25.0	
Chlorobenzene	1.019	1.094	0.500	-7.4	25.0	
Ethylbenzene	0.523	0.558	0.100	-6.7	25.0	
Styrene	0.981	1.075	0.300	-9.6	25.0	
Xylene (total)	0.633	0.692	0.300	-9.3	25.0	
Toluene-d8-	1.169	1.204		-3.0		
Bromofluorobenzene	0.808	0.834	0.200	-3.2	25.0	
1,2-Dichloroethane-d4	1.604	1.629		-1.6		

All other compounds must meet a minimum RRF of 0.010.

- 6A
VOLATILE ORGANICS INITIAL CALIBRATION DATA

Lab Name: COMPUCHEM RTP

Contract: 68D10083

ORIGINAL
(R90)

Lab Code: COMPU

Case No.: 18347

SAS No.: _____

SDG No.: CKY14

Instrument ID: F50056

Calibration Date(s): 06/25/92

06/25/92

Heated Purge: (Y/N): N

Calibration Times: 0021

0351

GC Column: DB624 ID: 0.530(mm)

LAB FILE ID:	RRF10 = <u>CU920625C56</u>	RRF20 = <u>CY920625C56</u>
RRF50=	<u>CS920625C56</u>	RRF100= <u>CW920625C56</u>
RRF200=	<u>CT920625C56</u>	

COMPOUND	RRF10	RRF20	RRF50	RRF100	RRF200	RRF	% RSD
Chloromethane	0.232	0.164	0.172	0.144	0.161	0.175	19.2
Bromomethane	* 1.621	1.257	1.156	1.215	1.092	1.268	16.3*
Vinyl Chloride	* 1.138	0.857	0.790	0.847	0.837	0.894	15.5*
Chloroethane	1.117	0.917	0.829	0.828	0.785	0.895	14.9
Methylene Chloride	2.523	1.777	1.806	1.566	1.370	1.808	24.2
Acetone	0.819	0.413	0.258	0.221	0.214	0.385	66.4
Carbon Disulfide	4.747	3.885	4.302	4.099	3.769	4.160	9.3
1,1-Dichloroethene	* 1.711	1.384	1.578	1.383	1.250	1.461	12.5*
1,1-Dichloroethane	* 3.684	3.084	3.377	3.159	2.955	3.252	8.8*
1,2-Dichloroethene (total)	1.857	1.520	1.658	1.562	1.437	1.607	10.0
Chlroform	* 3.636	3.032	3.287	3.074	2.876	3.181	9.2*
1,2-Dichloroethane	* 2.015	1.690	1.894	1.778	1.684	1.812	7.8*
2-Butanone	0.565	0.331	0.377	0.394	0.380	0.409	22.0
1,1-Trichloroethane	* 0.618	0.525	0.548	0.518	0.487	0.539	9.1*
Carbon Tetrachloride	* 0.520	0.450	0.496	0.484	0.473	0.485	5.4*
Bromodichloromethane	* 0.651	0.563	0.595	0.601	0.590	0.600	5.3*
1,2-Dichloropropane	0.492	0.425	0.453	0.437	0.416	0.445	6.7
cis-1,3-Dichloropropene	* 0.670	0.613	0.651	0.596	0.623	0.631	4.7*
Trichloroethene	* 0.525	0.446	0.473	0.449	0.427	0.464	8.2*
Dibromochloromethane	* 0.450	0.406	0.484	0.467	0.454	0.452	6.4*
1,1,2-Trichloroethane	* 0.440	0.340	0.364	0.337	0.304	0.357	14.3*
Benzene	* 1.214	1.003	1.062	0.995	0.918	1.038	10.7*
Trans-1,3-Dichloropropene	* 0.426	0.369	0.422	0.401	0.380	0.400	6.3*
Bromoform	* 0.311	0.276	0.325	0.319	0.319	0.310	6.3*
4-Methyl-2-Pentanone	0.381	0.356	0.324	0.331	0.361	0.351	6.6
2-Hexanone	0.259	0.155	0.174	0.193	0.183	0.193	20.5
Tetrachloroethene	* 0.572	0.484	0.506	0.471	0.441	0.495	9.9*
1,1,2,2-Tetrachloroethane	* 0.582	0.494	0.538	0.495	0.478	0.517	8.2*
Toluene	* 1.593	1.171	1.434	1.313	1.126	1.327	14.4*
Chlorobenzene	* 1.189	1.012	1.075	1.005	0.925	1.041	9.4*
Ethylbenzene	* 0.593	0.523	0.557	0.523	0.479	0.535	8.0*
Styrene	* 1.149	0.970	1.039	0.974	0.907	1.008	9.1*
Xylene (total)	* 0.731	0.632	0.671	0.618	0.564	0.643	9.7*
Toluene-d8	-	1.457	1.100	1.218	1.143	1.052	1.194
Bromofluorobenzene	* 1.015	0.889	0.873	0.847	0.805	0.886	8.9*
1,2-Dichloroethane-d4	2.048	1.706	1.706	1.707	1.690	1.771	8.7

* Compounds with required minimum RRF and maximum %RSD values.

All other compounds must meet a minimum RRF of 0.010.

7A
VOLATILE CONTINUING CALIBRATION CHECK

ORIGINAL
(Red)

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 18347 SAS No.: SDG No.: CKY14

Instrument ID: F50056 Calibration date: 06/25/92 Time: 0912

Lab File ID: CT920625A56 Init. Calib. Date(s): 06/25/92 06/25/92

Heated Purge: (Y/N) N Init. Calib. Times: 0021 0351

GC Column: DB624 ID: 0.530(mm)

COMPOUND	RRF	RRF50	MIN RRF	%D	%D	Samples
Chloromethane	0.175	0.185	-5.7			CKY14
Bromomethane	1.268	1.140	0.100	10.1	25.0	CKY17
Vinyl Chloride	0.894	0.819	0.100	8.4	25.0	CKY19
Chloroethane	0.895	0.809		9.6		CKY20
Methylene Chloride	1.808	1.805		0.2		CKY20
Acetone	0.385	0.197		48.8		CKY22
Carbon Disulfide	4.150	4.164		-0.1		CKY23
1,1-Dichloroethene	1.461	1.500	0.100	-2.7	25.0	
1,1-Dichloroethane	3.252	3.264	0.200	-0.4	25.0	CKY24
1,2-Dichloroethene (total)	1.607	1.644		-2.3		CKY25
Chloroform	3.181	3.298	0.200	-3.7	25.0	CKY18MS
1,2-Dichloroethane	1.812	1.699	0.100	6.2	25.0	CKY18MS
2-Butanone	0.409	0.247		39.6		CKY18MS
1,1,1-Trichloroethane	0.539	0.526	0.100	2.4	25.0	
Carbon Tetrachloride	0.485	0.483	0.100	0.4	25.0	
Bromodichloromethane	0.600	0.547	0.200	8.8	25.0	
1,2-Dichloropropane	0.445	0.421		5.4		
cis-1,3-Dichloropropene	0.631	0.583	0.200	7.6	25.0	
Trichloroethene	0.464	0.458	0.300	1.3	25.0	
Dibromochloromethane	0.452	0.454	0.100	-0.4	25.0	
1,1,2-Trichloroethane	0.357	0.327	0.100	8.4	25.0	
Benzene	1.038	1.079	0.500	-4.0	25.0	
Trans-1,3-Dichloropropene	0.400	0.369	0.100	7.8	25.0	
Bromoform	0.310	0.300	0.100	3.2	25.0	
4-Methyl-2-Pentanone	0.351	0.301		14.2		
2-Hexanone	0.193	0.154		20.2		
Tetrachloroethene	0.495	0.499	0.200	-0.8	25.0	
1,1,2,2-Tetrachloroethane	0.517	0.518	0.500	-0.2	25.0	
Toluene	1.327	1.411	0.400	-6.3	25.0	
Chlorobenzene	1.041	1.072	0.500	-3.0	25.0	
Ethylbenzene	0.535	0.553	0.100	-3.4	25.0	
Styrene	1.008	1.037	0.300	-2.9	25.0	
Xylene (total)	0.643	0.667	0.300	-3.7	25.0	
Toluene-d8-	1.194	1.248		-4.5		
Bromofluorobenzene	0.886	0.887	0.200	-0.1	25.0	
1,2-Dichloroethane-d4	1.771	1.657		6.4		

All other compounds must meet a minimum RRF of 0.010.

7A
VOLATILE CONTINUING CALIBRATION CHECK

ORIGINAL
(Red)

Lab Name: COMPUCHEM, RTP Contract: 68D10083

Lab Code: COMPU Case No.: 18347 SAS No.: SDG No.: CKY14

Instrument ID: F50056 Calibration date: 06/25/92 Time: 2225

Lab File ID: CS920625B56 Init. Calib. Date(s): 06/25/92 06/25/92

Heated Purge: (Y/N) N Init. Calib. Times: 0021 0351

GC Column: DB624 ID: 0.530 (mm)

COMPOUND	RRF	RRF50	MIN RRF	%D	MAX %D	Samp ¹⁴
Chloromethane	0.175	0.240	-37.1			CKY18
Bromomethane	1.268	1.335	0.100	-5.3	25.0	CKY26
Vinyl Chloride	0.894	1.053	0.100	-17.8	25.0	
Chloroethane	0.895	0.874		2.3		
Methylene Chloride	1.808	1.871		-3.5		
Acetone	0.385	0.249	35.3			
Carbon Disulfide	4.160	4.367		-5.0		
1,1-Dichloroethene	1.461	1.572	0.100	-7.6	25.0	
1,1-Dichloroethane	3.252	3.405	0.200	-4.7	25.0	
1,2-Dichloroethene (total)	1.607	1.702		-5.9		
Chloroform	3.181	3.390	0.200	-6.6	25.0	
1,2-Dichloroethane	1.812	1.898	0.100	-4.7	25.0	
2-Butanone	0.409	0.360		12.0		
1,1,1-Trichloroethane	0.539	0.545	0.100	-1.1	25.0	
Carbon Tetrachloride	0.485	0.490	0.100	-1.0	25.0	
Bromodichloromethane	0.600	0.613	0.200	-2.2	25.0	
1,2-Dichloropropane	0.445	0.466		-4.7		
cis-1,3-Dichloropropene	0.631	0.677	0.200	-7.3	25.0	
Trichloroethene	0.464	0.477	0.300	-2.8	25.0	
Dibromochloromethane	0.452	0.463	0.100	-2.4	25.0	
1,1,2-Trichloroethane	0.357	0.363	0.100	-1.7	25.0	
Benzene	1.038	1.095	0.500	-5.5	25.0	
Trans-1,3-Dichloropropene	0.400	0.402	0.100	-0.5	25.0	
Bromoform	0.310	0.294	0.100	5.2	25.0	
4-Methyl-2-Pentanone	0.351	0.360		-2.6		
2-Hexanone	0.193	0.158		18.1		
Tetrachloroethene	0.495	0.496	0.200	-0.2	25.0	
1,1,2,2-Tetrachloroethane	0.517	0.513	0.500	0.8	25.0	
Toluene	1.327	1.234	0.400	7.0	25.0	
Chlorobenzene	1.041	1.079	0.500	-3.6	25.0	
Ethylbenzene	0.535	0.554	0.100	-3.6	25.0	
Styrene	1.008	1.049	0.300	-4.1	25.0	
Xylene (total)	0.643	0.663	0.300	-3.1	25.0	
Toluene-d8	1.194	1.112		6.9		
Bromofluorobenzene	0.886	0.883	0.200	0.3	25.0	
1,2-Dichloroethane-d4	1.771	1.787		-0.9		

All other compounds must meet a minimum RRF of 0.010.

6B
SEMIVOLATILE ORGANICS INITIAL CALIBRATION DATA

Lab Name: COMPUCHEM, RTP Contract: 68D10083
 Lab Code: COMPU Case No.: 18347 SAS No.: SDG No.: CKY28
 Instrument ID: F50057 Calibration Date(s): 05/04/92 05/04/92
 Calibration Times: 1831 2137

ORIGINAL
(Red)

LAB FILE ID:	RRF20 = <u>HA920504B57</u>	RRF50 = <u>HB920504B57</u>	RRF80	RRF120	RRF160	RRF	% RSD	
	<u>RRF80 = HC920504B57</u>	<u>RRF120= HD920504B57</u>		<u>RRF160= HE920504B57</u>				
Phenol	*	1.899	1.961	1.821	1.800	1.766	1.849	4.3*
bis(2-Chloroethyl)Ether	*	1.680	1.684	1.570	1.543	1.479	1.591	5.6*
2-Chlorophenol	*	1.420	1.439	1.336	1.300	1.284	1.356	5.2*
1,3-Dichlorobenzene	*	1.531	1.493	1.372	1.323	1.236	1.391	8.7*
1,4-Dichlorobenzene	*	1.620	1.568	1.439	1.388	1.315	1.466	8.6*
1,2-Dichlorobenzene	*	1.480	1.458	1.354	1.312	1.284	1.378	6.3*
2-Methylphenol	*	1.299	1.386	1.340	1.305	1.273	1.321	3.3*
2,2'-Oxybis(1-Chloropropane)	2.625	2.854	2.792	2.857	2.843	2.794	3.5	
4-Methylphenol	*	1.322	1.422	1.323	1.325	1.308	1.340	3.5*
N-Nitroso-Di-n-Propylamine	*	1.489	1.553	1.482	1.518	1.518	1.512	1.9*
Hexachloroethane	*	0.740	0.765	0.714	0.707	0.699	0.725	3.7*
Nitrobenzene	*	0.500	0.494	0.473	0.457	0.449	0.475	4.7*
Isophorone	*	0.889	0.884	0.847	0.847	0.830	0.859	3.0*
2-Nitrophenol	*	0.163	0.184	0.176	0.176	0.172	0.174	4.4*
2,4-Dimethylphenol	*	0.339	0.311	0.319	0.319	0.311	0.320	3.6*
2-(2-Chloroethoxy)Methane	*	0.531	0.551	0.524	0.514	0.499	0.524	3.7*
2,4-Dichlorophenol	*	0.247	0.266	0.261	0.255	0.245	0.255	3.5*
1,2,4-Trichlorobenzene	*	0.301	0.295	0.289	0.268	0.254	0.281	7.0*
Naphthalene	*	1.044	0.994	0.921	0.873	0.822	0.931	9.6*
4-Chloroaniline		0.274	0.176	0.367	0.395	0.367	0.316	28.7
Hexachlorobutadiene		0.163	0.157	0.154	0.142	0.135	0.150	7.6
4-Chloro-3-Methylphenol	*	0.313	0.335	0.325	0.326	0.313	0.322	2.9*
2-Methylnaphthalene	*	0.628	0.629	0.594	0.550	0.519	0.584	8.3*
Hexachlorocyclopentadiene		0.348	0.353	0.359	0.346	0.321	0.345	4.2
2,4,6-Trichlorophenol	*	0.337	0.346	0.341	0.339	0.336	0.340	1.2*
2,4,5-Trichlorophenol	*	0.348	0.371	0.370	0.365	0.337	0.358	4.2*
2-Chloronaphthalene	*	1.104	1.120	1.076	1.016	0.965	1.056	6.1*
2-Nitroaniline		0.488	0.551	0.536	0.548	0.548	0.534	5.0
Dimethyl Phthalate		1.245	1.282	1.261	1.235	1.203	1.245	2.4
Acenaphthylene	*	1.779	1.774	1.716	1.621	1.545	1.687	6.0*
2,6-Dinitrotoluene	*	0.280	0.307	0.309	0.315	0.315	0.305	4.8*
3-Nitroaniline		0.207	0.260	0.264	0.299	0.308	0.268	14.9
Acenaphthene	*	1.102	1.104	1.047	0.981	0.913	1.029	8.0*
2,4-Dinitrophenol		0.107	0.163	0.166	0.172	0.165	0.155	17.3
4-Nitrophenol		0.128	0.167	0.162	0.167	0.162	0.157	10.5
Dibenzofuran		1.571	1.565	1.527	1.491	1.416	1.514	4.2*
2,4-Dinitrotoluene	*	0.365	0.408	0.411	0.419	0.407	0.402	5.3*

* Compounds with required minimum RRF and maximum %RSD values.
 All other compounds must meet a minimum RRF of 0.010.

6C
SEMIVOLATILE ORGANICS INITIAL CALIBRATION DATA

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

ORIGINAL
(Recd)

Lab Code: COMPU

Case No.: 18347

SAS No.: _____

SDG No.: CKY28

Instrument ID: F50057

Calibration Date(s): 05/04/92 05/04/92

Calibration Times: 1831 2137

LAB FILE ID:	RRF20 = HA920504B57	RRF50 = HB920504B57
RRF80 = HC920504B57	RRF120= HD920504B57	RRF160= HE920504B57

COMPOUND	RRF20	RRF50	RRF80	RRF120	RRF160	RRF	% RSD
Diethylphthalate	1.292	1.387	1.348	1.333	1.279	1.328	3.3
4-Chlorophenyl-phenylether*	0.607	0.623	0.612	0.580	0.541	0.593	5.5*
Fluorene	* 1.227	1.257	1.202	1.166	1.096	1.190	5.2*
4-Nitroaniline	0.190	0.236	0.237	0.282	0.302	0.249	17.6
4,6-Dinitro-2-Methylphenol	0.109	0.138	0.143	0.156	0.162	0.142	14.5
N-Nitrosodiphenylamine (1)	0.505	0.411	0.496	0.519	0.545	0.495	10.2
4-Bromophenyl-phenylether	* 0.216	0.217	0.218	0.213	0.218	0.216	1.0*
Hexachlorobenzene	* 0.261	0.248	0.241	0.232	0.236	0.244	4.7*
Pentachlorophenol	* 0.111	0.134	0.135	0.135	0.133	0.130	8.0*
Phenanthrene	* 1.031	1.045	1.013	0.960	0.921	0.994	5.2*
Anthracene	* 0.970	0.953	0.959	0.908	0.876	0.933	4.3*
Carbazole	0.630	0.581	0.494	0.525	0.580	0.562	9.5
Di-n-Butylphthalate	1.298	1.450	1.377	1.339	1.259	1.345	5.5
Fluoranthene	* 0.932	0.955	0.918	0.872	0.860	0.907	4.4*
Pyrene	* 1.449	1.509	1.383	1.339	1.250	1.386	7.2*
Butylbenzylphthalate	0.704	0.809	0.758	0.779	0.779	0.766	5.1
3,3'-Dichlorobenzidine	0.165	0.134	0.167	0.178	0.179	0.165	11.0
Benzo(a)Anthracene	* 1.042	1.089	1.046	1.020	1.016	1.043	2.8*
Chrysene	* 0.986	0.960	0.883	0.897	0.826	0.910	7.0*
bis(2-Ethylhexyl)Phthalate	0.997	1.124	1.051	1.097	1.088	1.071	4.6
Di-n-Octyl Phthalate	1.788	1.949	1.812	1.854	1.925	1.866	3.7
Benzo(b)Fluoranthene	* 1.059	1.100	1.036	1.073	1.123	1.078	3.2*
Benzo(k)Fluoranthene	* 0.973	1.027	0.935	0.842	0.807	0.917	9.9*
Benzo(a)Pyrene	* 0.897	0.928	0.928	0.952	0.942	0.929	2.2*
Indeno(1,2,3-cd)Pyrene	* 0.785	0.828	0.799	0.796	0.802	0.802	2.0*
Dibenz(a,h)Anthracene	* 0.792	0.703	0.688	0.683	0.661	0.705	7.2*
Benzo(g,h,i)Perylene	* 0.862	0.737	0.727	0.678	0.669	0.735	10.5*
Nitrobenzene-d5	* 0.464	0.474	0.440	0.438	0.426	0.448	4.4*
2-Fluorobiphenyl	* 1.268	1.211	1.149	1.080	1.012	1.144	8.9*
Terphenyl-d14	* 0.988	1.031	0.922	0.893	0.844	0.936	8.0*
Phenol-d5	* 1.792	1.804	1.710	1.654	1.604	1.713	5.0*
2-Fluorophenol	* 1.270	1.292	1.169	1.172	1.122	1.205	6.0*
2,4,6-Tribromophenol	0.185	0.189	0.184	0.180	0.177	0.183	2.5
2-Chlorophenol-d4	* 1.340	1.400	1.283	1.250	1.196	1.294	6.1*
1,2-Dichlorobenzene-d4	* 0.886	0.865	0.811	0.770	0.735	0.813	7.8*

(1) Cannot be separated from Diphenylamine

* Compounds with required minimum RRF and maximum %RSD values.

All other compounds must meet a minimum RRF of 0.010.

7B
SEMIVOLATILE CONTINUING CALIBRATION CHECK

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

ORIGINAL
(Red)

Lab Code: COMPU

Case No.: 18347

SAS No.: _____

SDG No.: CKY28

Instrument ID: F50057

Calibration date: 07/08/92 Time: 2010

Lab File ID: HG920708B57

Init. Calib. Date(s): 05/04/92 05/04/92

Init. Calib. Times: 1831 2137

COMPOUND	RRF	RRF50	MIN RRF	%D	MAX %D	Sample
Phenol	1.849	1.969	0.800	-6.5	25.0	
bis(2-Chloroethyl)Ether	1.591	1.689	0.700	-6.2	25.0	
2-Chlorophenol	1.356	1.454	0.800	-7.2	25.0	
1,3-Dichlorobenzene	1.391	1.567	0.600	-12.6	25.0	
1,4-Dichlorobenzene	1.466	1.594	0.500	-8.7	25.0	
1,2-Dichlorobenzene	1.378	1.439	0.400	-4.4	25.0	
2-Methylphenol	1.321	1.380	0.700	-4.5	25.0	
2,2'-Oxybis(1-Chloropropane)	2.794	2.873		-2.8		
4-Methylphenol	1.340	1.447	0.600	-8.0	25.0	
N-Nitroso-Di-n-Propylamine	1.512	1.491	0.500	1.4	25.0	
Hexachloroethane	0.725	0.803	0.300	-10.8	25.0	
Nitrobenzene	0.475	0.507	0.200	-6.7	25.0	
Isophorone	0.859	0.928	0.400	-8.0	25.0	
2-Nitrophenol	0.174	0.193	0.100	-10.9	25.0	
2,4-Dimethylphenol	0.320	0.345	0.200	-7.8	25.0	
bis(2-Chloroethoxy)Methane	0.524	0.515	0.300	1.7	25.0	
2,4-Dichlorophenol	0.255	0.288	0.200	-12.9	25.0	
1,2,4-Trichlorobenzene	0.281	0.299	0.200	-6.4	25.0	
Naphthalene	0.931	1.016	0.700	-9.1	25.0	
4-Chloroaniline	0.316	0.384		-21.5		
Hexachlorobutadiene	0.150	0.160		-6.7		
4-Chloro-3-Methylphenol	0.322	0.361	0.200	-12.1	25.0	
2-Methylnaphthalene	0.584	0.659	0.400	-12.8	25.0	
Hexachlorocyclopentadiene	0.345	0.367		-6.4		
2,4,6-Trichlorophenol	0.340	0.353	0.200	-3.8	25.0	
2,4,5-Trichlorophenol	0.358	0.364	0.200	-1.7	25.0	
2-Chloronaphthalene	1.056	1.169	0.800	-10.7	25.0	
2-Nitroaniline	0.534	0.538		-0.7		
Dimethyl Phthalate	1.245	1.399		-12.4		
Acenaphthylene	1.687	1.831	1.300	-8.5	25.0	
2,6-Dinitrotoluene	0.305	0.350	0.200	-14.8	25.0	
3-Nitroaniline	0.268	0.302		-12.7		
Acenaphthene	1.029	1.064	0.800	-3.4	25.0	
2,4-Dinitrophenol	0.155	0.165		-6.5		
4-Nitrophenol	0.157	0.180		-14.6		
Dibenzofuran	1.514	1.630	0.800	-7.7	25.0	
2,4-Dinitrotoluene	0.402	0.439	0.200	-9.2	25.0	

(1) Cannot be separated from Diphenylamine

All other compounds must meet a minimum RRF of 0.010.

7C
SEMICVOLATILE CONTINUING CALIBRATION CHECK

ORIGINAL
(Rev)

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 18347 SAS No.: SDG No.: CKY28

Instrument ID: F50057 Calibration date: 07/08/92 Time: 2010

Lab File ID: HG920708357 Init. Calib. Date(s): 05/04/92 05/04/92

Init. Calib. Times: 1831 2137

COMPOUND	RRF	RRF50	MIN RRF	%D	MAX %D
Diethylphthalate	1.328	1.486	-11.9		
4-Chlorophenyl-phenylether	0.593	0.647	0.400	-9.1	25.0
Fluorene	1.190	1.214	0.900	-2.0	25.0
4-Nitroaniline	0.249	0.252		-1.2	
4,6-Dinitro-2-Methylphenol	0.142	0.144		-1.4	
N-Nitrosodiphenylamine (1)	0.495	0.544		-9.9	
4-Bromophenyl-phenylether	0.216	0.239	0.100	-10.6	25.0
Hexachlorobenzene	0.244	0.251	0.100	-2.9	25.0
Pentachlorophenol	0.130	0.150	0.050	-15.4	25.0
Phenanthrene	0.994	1.059	0.700	-6.5	25.0
Anthracene	0.933	1.056	0.700	-13.2	25.0
Carbazole	0.562	0.670		-19.2	
Di-n-Butylphthalate	1.345	1.534		-14.0	
Fluoranthene	0.907	1.147	0.600	-26.5	25.0
Pyrene	1.386	1.483	0.600	-7.0	25.0
Butylbenzylphthalate	0.766	0.824		-7.6	
3,3'-Dichlorobenzidine	0.165	0.224		-35.8	
Benzo(a)Anthracene	1.043	1.150	0.800	-10.3	25.0
Chrysene	0.910	1.072	0.700	-17.8	25.0
bis(2-Ethylhexyl)Phthalate	1.071	1.110		-3.6	
Di-n-Octyl Phthalate	1.866	1.855		0.6	
Benzo(b)Fluoranthene	1.078	1.166	0.700	-6.2	25.0
Benzo(k)Fluoranthene	0.917	1.122	0.700	-22.4	25.0
Benzo(a)Pyrene	0.929	1.045	0.700	-12.5	25.0
Indeno(1,2,3-cd)Pyrene	0.802	0.684	0.500	14.7	25.0
Dibenz(a,h)Anthracene	0.705	0.634	0.400	10.1	25.0
Benzo(g,h,i)Perylene	0.735	0.697	0.500	5.2	25.0
Nitrobenzene-d5	0.448	0.560	0.200	-25.0	25.0
2-Fluorobiphenyl	1.144	1.209	0.700	-5.7	25.0
Terphenyl-d14	0.936	1.006	0.500	-7.5	25.0
Phenol-d5	1.713	1.953	0.800	-14.0	25.0
2-Fluorophenol	1.205	1.542	0.600	-28.0	25.0
2,4,6-Tribromophenol	0.183	0.197		-7.6	
2-Chlorophenol-d4	1.294	1.296	0.800	-0.2	25.0
1,2-Dichlorobenzene-d4	0.813	0.918	0.400	-12.9	25.0

(1) Cannot be separated from Diphenylamine
All other compounds must meet a minimum RRF of 0.010.

7B
SEMIVOLATILE CONTINUING CALIBRATION CHECK

ORIGINAL
(Red)

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 18347 SAS No.: _____ SDG No.: CKY28

Instrument ID: F50057 Calibration date: 07/09/92 Time: 1030

Lab File ID: HG920709A57 Init. Calib. Date(s): 05/04/92 05/04/92

Init. Calib. Times: 1831 2137

COMPOUND	RRF	RRF50	MIN RRF	%D	MAX %D	Sample
Phenol	1.849	2.010	0.800	-8.7	25.0	CKY28
bis(2-Chloroethyl)Ether	1.591	1.690	0.700	-6.2	25.0	CKY31
2-Chlorophenol	1.356	1.509	0.800	-11.3	25.0	CKY35
1,3-Dichlorobenzene	1.391	1.596	0.600	-14.7	25.0	CKY36
1,4-Dichlorobenzene	1.466	1.650	0.500	-12.6	25.0	CKY38
1,2-Dichlorobenzene	1.378	1.504	0.400	-9.1	25.0	
2-Methylphenol	1.321	1.371	0.700	-3.8	25.0	
2,2'-Oxybis(1-Chloropropane)	2.794	2.879		-3.0		
4-Methylphenol	1.340	1.424	0.600	-6.3	25.0	
N-Nitroso-Di-n-Propylamine	1.512	1.490	0.500	1.5	25.0	
Hexachloroethane	0.725	0.845	0.300	-16.6	25.0	
Nitrobenzene	0.475	0.552	0.200	-16.2	25.0	
Isophorone	0.859	0.956	0.400	-11.3	25.0	
2-Nitrophenol	0.174	0.197	0.100	-13.2	25.0	
2,4-Dimethylphenol	0.320	0.359	0.200	-12.2	25.0	
bis(2-Chloroethoxy)Methane	0.524	0.517	0.300	1.3	25.0	
2,4-Dichlorophenol	0.255	0.292	0.200	-14.5	25.0	
1,2,4-Trichlorobenzene	0.281	0.300	0.200	-6.8	25.0	
Naphthalene	0.931	1.032	0.700	-10.8	25.0	
4-Chloroaniline	0.316	0.367		-16.1		
Hexachlorobutadiene	0.150	0.162		-8.0		
4-Chloro-3-Methylphenol	0.322	0.368	0.200	-14.3	25.0	
2-Methylnaphthalene	0.584	0.676	0.400	-15.8	25.0	
Hexachlorocyclopentadiene	0.345	0.364		-5.5		
2,4,6-Trichlorophenol	0.340	0.352	0.200	-3.5	25.0	
2,4,5-Trichlorophenol	0.358	0.364	0.200	-1.7	25.0	
2-Chloronaphthalene	1.056	1.193	0.800	-13.0	25.0	
2-Nitroaniline	0.534	0.551		-3.2		
Dimethyl Phthalate	1.245	1.392		-11.8		
Acenaphthylene	1.687	1.864	1.300	-10.5	25.0	
2,6-Dinitrotoluene	0.305	0.342	0.200	-12.1	25.0	
3-Nitroaniline	0.268	0.282		-5.2		
Acenaphthene	1.029	1.113	0.800	-8.2	25.0	
2,4-Dinitrophenol	0.155	0.150		3.2		
4-Nitrophenol	0.157	0.185		-17.8		
Dibenzofuran	1.514	1.783	0.800	-17.8	25.0	
2,4-Dinitrotoluene	0.402	0.428	0.200	-6.5	25.0	

(1) Cannot be separated from Diphenylamine

All other compounds must meet a minimum RRF of 0.010.

7C
SEMIVOLATILE CONTINUING CALIBRATION CHECK

ORIGINAL
(Rev)

Lab Name: COMPUCHEM RTP

Contract: 68D10083

Lab Code: COMPU

Case No.: 18347

SAS No.: _____

SDG No.: CKY28

Instrument ID: F50057

Calibration date: 07/09/92 Time: 1030

Lab File ID: HG920709A57

Init. Calib. Date(s): 05/04/92 05/04/92

Init. Calib. Times: 1831 2137

COMPOUND	RRF	RRF50	MIN RRF	%D	MAX %D
Diethylphthalate	1.328	1.709		-28.7	
4-Chlorophenyl-phenylether	0.593	0.692	0.400	-16.7	25.0
Fluorene	1.190	1.340	0.900	-12.6	25.0
4-Nitroaniline	0.249	0.271		-8.8	
4,6-Dinitro-2-Methylphenol	0.142	0.145		-2.1	
N-Nitrosodiphenylamine (1)	0.495	0.563		-13.7	
4-Bromophenyl-phenylether	0.216	0.238	0.100	-10.2	25.0
Hexachlorobenzene	0.244	0.258	0.100	-5.7	25.0
Pentachlorophenol	0.130	0.152	0.050	-16.9	25.0
Phenanthrene	0.994	1.076	0.700	-8.2	25.0
Anthracene	0.933	1.053	0.700	-12.9	25.0
Carbazole	0.562	0.658		-17.1	
Di-n-Butylphthalate	1.345	1.588		-18.1	
Fluoranthene	0.907	1.134	0.600	-25.0	25.0
Pyrene	1.386	1.423	0.600	-2.7	25.0
Butylbenzylphthalate	0.766	0.793		-3.5	
3,3'-Dichlorobenzidine	0.165	0.264		-60.0	
Benzo(a)Anthracene	1.043	1.120	0.800	-7.4	25.0
Chrysene	0.910	1.034	0.700	-13.6	25.0
bis(2-Ethylhexyl)Phthalate	1.071	1.006		6.1	
Di-n-Octyl Phthalate	1.866	1.457		21.9	
Benzo(b)Fluoranthene	1.078	1.145	0.700	-5.2	25.0
Benzo(k)Fluoranthene	0.917	1.011	0.700	-10.2	25.0
Benzo(a)Pyrene	0.929	1.033	0.700	-11.2	25.0
Indeno(1,2,3-cd)Pyrene	0.802	0.850	0.500	-6.0	25.0
Dibenz(a,h)Anthracene	0.705	0.740	0.400	-5.0	25.0
Benzo(g,h,i)Perylene	0.735	0.808	0.500	-9.9	25.0
Nitrobenzene-d5	0.448	0.618	0.200	-38.0	25.0
2-Fluorobiphenyl	1.144	1.238	0.700	-8.2	25.0
Terphenyl-d14	0.936	0.965	0.500	-3.1	25.0
Phenol-d5	1.713	1.984	0.800	-15.8	25.0
2-Fluorophenol	1.205	1.534	0.600	-27.3	25.0
2,4,6-Tribromophenol	0.183	0.199		-8.7	
2-Chlorophenol-d4	1.294	1.280	0.800	1.1	25.0
1,2-Dichlorobenzene-d4	0.813	0.942	0.400	-15.9	25.0

(1) Cannot be separated from Diphenylamine
All other compounds must meet a minimum RRF of 0.010.

6B
SEMIVOLATILE ORGANICS INITIAL CALIBRATION DATA

ORIGINAL
(Red)

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU

Case No.: 18347

SAS No.: _____

SDG No.: CKY28

Instrument ID: OWA15

Calibration Date(s): 06/11/92 06/11/92

Calibration Times: 1923 2357

LAB FILE ID:	RRF20 = <u>HI920611B15</u>	RRF50 = <u>HG920611B15</u>
RRF30 = <u>HK920611B15</u>	RRF120= <u>HJ920611B15</u>	RRF160= <u>HH920611B15</u>

COMPOUND	RRF20	RRF50	RRF80	RRF120	RRF160	RRF	% RSD
Phenol	* 1.533	1.555	1.566	1.610	1.493	1.551	2.8*
bis(2-Chloroethyl)Ether	* 1.374	1.427	1.421	1.451	1.492	1.433	3.0*
2-Chlorophenol	* 1.324	1.379	1.380	1.205	1.383	1.334	5.7*
1,3-Dichlorobenzene	* 1.454	1.443	1.422	1.502	1.445	1.453	2.0*
1,4-Dichlorobenzene	* 1.566	1.547	1.508	1.565	1.559	1.549	1.6*
1,2-Dichlorobenzene	* 1.430	1.396	1.402	1.464	1.420	1.422	1.9*
2-Methylphenol	* 1.073	1.119	1.189	1.184	1.171	1.147	4.4*
2,2'-Oxybis(1-Chloropropane)	1.524	1.640	1.720	1.749	1.901	1.707	8.2
4-Methylphenol	* 1.088	1.191	1.200	1.247	1.238	1.193	5.3*
N-Nitroso-Di-n-Propylamine	* 1.016	1.116	1.126	1.168	1.204	1.126	6.3*
Hexachloroethane	* 0.873	0.842	0.882	0.903	0.852	0.870	2.8*
Nitrobenzene	* 0.418	0.467	0.463	0.469	0.500	0.463	6.3*
Isophorone	* 0.745	0.872	0.860	0.884	0.930	0.858	8.0*
2-Nitrophenol	* 0.171	0.197	0.194	0.191	0.199	0.190	5.9*
2,4-Dimethylphenol	* 0.359	0.354	0.394	0.382	0.398	0.377	5.3*
bis(2-Chloroethoxy)Methane	* 0.487	0.522	0.530	0.539	0.541	0.524	4.2*
2,4-Dichlorophenol	* 0.260	0.299	0.297	0.305	0.304	0.293	6.4*
1,2,4-Trichlorobenzene	* 0.314	0.326	0.323	0.329	0.335	0.325	2.4*
Naphthalene	* 1.033	1.061	1.089	1.101	1.048	1.066	2.6*
4-Chloroaniline	0.324	0.100	0.346	0.311	0.360	0.288	37.1
Hexachlorobutadiene	0.188	0.187	0.193	0.193	0.199	0.192	2.5
4-Chloro-3-Methylphenol	* 0.265	0.317	0.310	0.311	0.338	0.308	8.7*
2-Methylnaphthalene	* 0.623	0.663	0.672	0.677	0.693	0.666	3.9*
Hexachlorocyclopentadiene	0.343	0.363	0.418	0.437	0.460	0.404	12.3
2,4,6-Trichlorophenol	* 0.320	0.374	0.377	0.380	0.424	0.375	9.8*
2,4,5-Trichlorophenol	* 0.318	0.354	0.361	0.360	0.361	0.351	5.3*
2-Chloronaphthalene	* 1.187	1.208	1.280	1.273	1.285	1.247	3.7*
2-Nitroaniline	0.238	0.339	0.352	0.358	0.412	0.340	18.6
Dimethyl Phthalate	1.185	1.295	1.316	1.269	1.340	1.281	4.7
Acenaphthylene	* 1.782	1.906	2.017	1.988	2.049	1.948	5.5*
2,6-Dinitrotoluene	* 0.205	0.262	0.261	0.260	0.285	0.255	11.6*
3-Nitroaniline	0.157	0.194	0.206	0.188	0.193	0.188	9.8
Acenaphthene	* 1.109	1.135	1.172	1.177	1.211	1.161	3.4*
2,4-Dinitrophenol	0.052	0.094	0.096	0.105	0.128	0.095	29.0
4-Nitrophenol	0.118	0.167	0.193	0.195	0.195	0.174	19.1
Dibenzofuran	* 1.483	1.548	1.547	1.548	1.624	1.550	3.2*
2,4-Dinitrotoluene	* 0.259	0.321	0.325	0.321	0.348	0.315	10.5*

(1) Cannot be separated from Diphenylamine

* Compounds with required minimum RRF and maximum %RSD values.

All other compounds must meet a minimum RRF of 0.010.

6C
SEMIVOLATILE ORGANICS INITIAL CALIBRATION DATA

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

ORIGINAL
(RSD)

Lab Code: COMPU

Case No.: 18347

SAS No.: _____

SDG No.: CKY28

Instrument ID: OWA15

Calibration Date(s): 06/11/92 06/11/92

Calibration Times: 1923 2357

LAB FILE ID:	RRF20 = HJ920611B15	RRF50 = HG920611B15
RRF80 = HK920611B15	RRF120= HJ920611B15	RRF160= HH920611B15

COMPOUND	RRF20	RRF50	RRF80	RRF120	RRF160	RRF	% RSD
Diethylphthalate	1.276	1.425	1.442	1.413	1.463	1.404	5.3
4-Chlorophenyl-phenylether*	0.490	0.525	0.539	0.549	0.584	0.537	6.4*
Fluorene	* 1.104	1.201	1.226	1.232	1.286	1.210	5.5*
4-Nitroaniline	0.131	0.160	0.174	0.162	0.155	0.156	10.1
4,6-Dinitro-2-Methylphenol	0.090	0.122	0.131	0.134	0.147	0.125	17.1
N-Nitrosodiphenylamine (1)	0.603	0.463	0.672	0.647	0.619	0.601	13.6
4-Bromophenyl-phenylether	* 0.200	0.216	0.218	0.217	0.219	0.214	3.7*
Hexachlorobenzene	* 0.281	0.279	0.280	0.271	0.285	0.279	1.8*
Pentachlorophenol	* 0.075	0.111	0.113	0.124	0.133	0.111	19.9*
Phenanthrene	* 1.237	1.223	1.302	1.271	1.308	1.268	3.0*
Anthracene	* 1.060	0.996	1.113	1.121	1.079	1.074	4.7*
Carbazole	0.741	0.760	0.862	0.820	0.753	0.787	6.6
Di-n-Butylphthalate	1.437	1.685	1.744	1.778	1.688	1.666	8.0
Fluoranthene	* 0.744	0.863	0.918	0.913	0.908	0.869	8.4*
Pyrene	* 2.037	2.283	2.118	2.138	2.148	2.145	4.1*
Butylbenzylphthalate	0.865	1.207	1.109	1.132	1.172	1.097	12.3
3,3'-Dichlorobenzidine	0.166	0.142	0.216	0.220	0.216	0.192	18.6
Benzo(a)Anthracene	* 1.135	1.216	1.219	1.219	1.266	1.211	3.9*
Chrysene	* 1.188	1.168	1.111	1.148	1.029	1.129	5.5*
bis(2-Ethylhexyl)Phthalate	1.057	1.509	1.422	1.542	1.549	1.416	14.6
Di-n-Octyl Phthalate	2.284	2.993	2.770	2.727	2.659	2.687	9.6
Benzo(b)Fluoranthene	* 1.190	1.376	1.218	1.451	1.401	1.327	8.7*
Benzo(k)Fluoranthene	* 1.273	1.161	1.226	0.915	0.850	1.085	17.6*
Benzo(a)Pyrene	* 0.952	1.037	1.081	1.090	1.029	1.038	5.3*
Indeno(1,2,3-cd)Pyrene	* 0.881	1.106	1.160	1.171	1.060	1.076	10.9*
Dibenz(a,h)Anthracene	* 0.669	0.898	0.958	0.961	0.907	0.879	13.7*
Benzo(g,h,i)Perylene	* 0.741	1.009	1.023	1.044	0.935	0.950	13.1*
Nitrobenzene-d5	* 0.393	0.470	0.467	0.469	0.508	0.461	9.1*
2-Fluorobiphenyl	* 1.417	1.401	1.481	1.488	1.495	1.456	3.0*
Terphenyl-d14	* 1.044	1.146	1.110	1.101	1.130	1.106	3.5*
Phenol-d5	* 1.466	1.585	1.569	1.573	1.641	1.567	4.0*
2-Fluorophenol	* 1.320	1.421	1.363	1.368	1.412	1.377	3.0*
2,4,6-Tribromophenol	0.085	0.116	0.112	0.117	0.132	0.112	15.3
2-Chlorophenol-d4	* 1.135	1.210	1.210	1.251	1.228	1.207	3.6*
1,2-Dichlorobenzene-d4	* 0.987	0.928	0.941	0.949	0.918	0.945	2.8*

(L) Cannot be separated from Diphenylamine

* Compounds with required minimum RRF and maximum %RSD values.

All other compounds must meet a minimum RRF of 0.010.

7B
SEMIVOLATILE CONTINUING CALIBRATION CHECK

ORIGINAL
(Red)

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 18347 SAS No.: _____ SDG No.: CKY28

Instrument ID: OWA15 Calibration date: 07/01/92 Time: 1718

Lab File ID: HG920701B15 Init. Calib. Date(s): 06/11/92 06/11/92

Init. Calib. Times: 1923 2357

COMPOUND	RRF	RRF50	MIN RRF	%D	MAX %D	Samples
Phenol	1.551	1.849	0.800	-19.2	25.0	CK430
bis(2-Chloroethyl)Ether	1.433	1.730	0.700	-20.7	25.0	CK432
2-Chlorophenol	1.334	1.431	0.800	-7.3	25.0	CK432MS
1,3-Dichlorobenzene	1.453	1.558	0.600	-7.2	25.0	CK432MSD
1,4-Dichlorobenzene	1.549	1.523	0.500	1.7	25.0	CK432MSD
1,2-Dichlorobenzene	1.422	1.382	0.400	2.8	25.0	CK433
2-Methylphenol	1.147	1.171	0.700	-2.1	25.0	CK433
2,2'-Oxybis(1-Chloropropane)	1.707	3.076		(-80.2)		CK437
4-Methylphenol	1.193	1.293	0.600	-8.4	25.0	
N-Nitroso-Di-n-Propylamine	1.126	1.375	0.500	-22.1	25.0	
Hexachloroethane	0.870	0.665	0.300	23.6	25.0	
Nitrobenzene	0.463	0.476	0.200	-2.8	25.0	
Isophorone	0.858	0.909	0.400	-5.9	25.0	
2-Nitrophenol	0.190	0.204	0.100	-7.4	25.0	
2,4-Dimethylphenol	0.377	0.301	0.200	20.2	25.0	
bis(2-Chloroethoxy)Methane	0.524	0.599	0.300	-14.3	25.0	
2,4-Dichlorophenol	0.293	0.301	0.200	-2.7	25.0	
1,2,4-Trichlorobenzene	0.325	0.337	0.200	-3.7	25.0	
Naphthalene	1.066	1.024	0.700	3.9	25.0	
4-Chloroaniline	0.288	0.256		11.1		
Hexachlorobutadiene	0.192	0.169		12.0		
4-Chloro-3-Methylphenol	0.308	0.299	0.200	2.9	25.0	
2-Methylnaphthalene	0.666	0.639	0.400	4.1	25.0	
Hexachlorocyclopentadiene	0.404	0.376		6.9		
2,4,6-Trichlorophenol	0.375	0.416	0.200	-10.9	25.0	
2,4,5-Trichlorophenol	0.351	0.431	0.200	-22.8	25.0	
2-Chloronaphthalene	1.247	1.241	0.800	0.5	25.0	
2-Nitroaniline	0.340	0.436		(-28.2)		
Dimethyl Phthalate	1.281	1.274		0.5		
Acenaphthylene	1.948	1.918	1.200	1.5	25.0	
2,6-Dinitrotoluene	0.255	0.281	0.200	-10.2	25.0	
3-Nitroaniline	0.188	0.259		(-37.8)		
Acenaphthene	1.161	1.156	0.800	0.4	25.0	
2,4-Dinitrophenol	0.095	0.136		(-43.2)		
4-Nitrophenol	0.174	0.085		(51.2)		
Dibenzofuran	1.550	1.557	0.800	-0.5	25.0	
2,4-Dinitrotoluene	0.315	0.345	0.200	-9.5	25.0	

All other compounds must meet a minimum RRF of 0.010.

7C
SEMIVOLATILE CONTINUING CALIBRATION CHECK

ORIGINAL
(Req)

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 18347 SAS No.: SDG No.: CKY28

Instrument ID: OWA15 Calibration date: 07/01/92 Time: 1718

Lab File ID: HG920701B15 Init. Calib. Date(s): 06/11/92 06/11/92

Init. Calib. Times: 1923 2357

COMPOUND	RRF	RRF50	MIN RRF	%D	MAX %D
Diethylphthalate	1.404	1.312		6.6	
4-Chlorophenyl-phenylether	0.537	0.612	0.400	-14.0	25.0
Fluorene	1.210	1.219	0.900	-0.7	25.0
4-Nitroaniline	0.156	0.241		(-54.5)	
4,6-Dinitro-2-Methylphenol	0.125	0.139		-11.2	
N-Nitrosodiphenylamine (1)	0.601	0.367		(38.9)	
4-Bromophenyl-phenylether	0.214	0.205	0.100	4.2	25.0
Hexachlorobenzene	0.279	0.252	0.100	9.7	25.0
Pentachlorophenol	0.111	0.123	0.050	-10.8	25.0
Phenanthrone	1.268	1.147	0.700	9.5	25.0
Anthracene	1.074	1.087	0.700	-1.2	25.0
Carbazole	0.787	0.876		-11.3	
Di-n-Butylphthalate	1.666	1.527		8.3	
Fluoranthene	0.869	0.986	0.600	-13.5	25.0
Pyrene	2.145	1.874	0.600	12.6	25.0
Butylbenzylphthalate	1.097	0.889		19.0	
3,3'-Dichlorobenzidine	0.192	0.174		9.4	
Benzo(a)Anthracene	1.211	1.199	0.800	1.0	25.0
Chrysene	1.129	1.039	0.700	8.0	25.0
bis(2-Ethylhexyl)Phthalate	1.416	1.174		17.1	
Di-n-Octyl Phthalate	2.687	2.130		20.7	
Benzo(b)Fluoranthene	1.327	1.133	0.700	14.6	25.0
Benzo(k)Fluoranthene	1.085	1.144	0.700	-5.4	25.0
Benzo(a)Pyrene	1.038	0.968	0.700	6.7	25.0
Indeno(1,2,3-cd)Pyrene	1.076	0.916	0.500	14.9	25.0
Dibenz(a,h)Anthracene	0.879	0.806	0.400	8.3	25.0
Benzo(g,h,i)Perylene	0.950	0.812	0.500	14.5	25.0
Nitrobenzene-d5	0.461	0.473	0.200	-2.6	25.0
2-Fluorobiphenyl	1.456	1.447	0.700	0.6	25.0
Terphenyl-d14	1.106	1.006	0.500	9.0	25.0
Phenol-d5	1.567	1.758	0.800	-12.2	25.0
2-Fluorophenol	1.377	1.376	0.600	0.1	25.0
2,4,6-Tribromophenol	0.112	0.132		-17.9	
2-Chlorophenol-d4	1.207	1.301	0.800	-7.8	25.0
1,2-Dichlorobenzene-d4	0.945	0.873	0.400	7.6	25.0

(1) Cannot be separated from Diphenylamine
All other compounds must meet a minimum RRF of 0.010.

✓

6B
SEMIVOLATILE ORGANICS INITIAL CALIBRATION DATA

Lab Name: COMPUCHEM RTP

Contract: 68D10083

ORIGINAL
(Rev)

Lab Code: COMPU

Case No.: 18347

SAS No.: _____

SDG No.: CKY28

Instrument ID: OWA05

Calibration Date(s): 06/11/92 06/12/92

Calibration Times: 2352 0249

<u>LAB FILE ID:</u>	<u>RRF20 = HI920612C05</u>	<u>RRF50 = HG920611B05</u>
<u>RRF80 = HK920612C05</u>	<u>RRF120= HJ920612C05</u>	<u>RRF160= HH920612C05</u>

COMPOUND	RRF20	RRF50	RRF80	RRF120	RRF160	RRF	% RSD
Phenol	* 1.822	1.846	1.846	1.822	1.740	1.815	2.4*
bis(2-Chloroethyl)Ether	* 1.608	1.626	1.593	1.557	1.526	1.582	2.5*
2-Chlorophenol	* 1.506	1.480	1.455	1.428	1.448	1.463	2.1*
1,3-Dichlorobenzene	* 1.590	1.499	1.482	1.486	1.501	1.512	2.9*
1,4-Dichlorobenzene	* 1.600	1.542	1.542	1.529	1.593	1.561	2.1*
1,2-Dichlorobenzene	* 1.437	1.362	1.355	1.387	1.461	1.400	3.3*
2-Methylphenol	* 1.399	1.434	1.377	1.370	1.325	1.381	2.9*
2,2'-Oxybis(1-Chloropropane)	2.201	2.145	2.178	2.098	1.960	2.116	4.5
4-Methylphenol	* 1.449	1.404	1.417	1.392	1.334	1.399	3.0*
N-Nitroso-Di-n-Propylamine	* 1.129	1.126	1.168	1.118	1.103	1.129	2.1*
Hexachloroethane	* 0.915	0.914	0.911	0.906	0.926	0.914	0.8*
Nitrobenzene	* 0.470	0.476	0.468	0.456	0.464	0.467	1.6*
Isophorone	* 0.870	0.889	0.874	0.857	0.894	0.877	1.7*
2-Nitrophenol	* 0.198	0.193	0.192	0.190	0.203	0.195	2.7*
2,4-Dimethylphenol	* 0.394	0.410	0.387	0.370	0.424	0.397	5.2*
(2-Chloroethoxy)Methane	* 0.561	0.549	0.538	0.520	0.520	0.538	3.3*
2,4-Dichlorophenol	* 0.316	0.312	0.300	0.301	0.312	0.308	2.3*
1,2,4-Trichlorobenzene	* 0.336	0.323	0.309	0.311	0.335	0.323	4.0*
Naphthalene	* 1.103	1.034	1.001	0.996	1.052	1.037	4.2*
4-Chloroaniline	0.455	0.438	0.430	0.360	0.405	0.418	8.8
Hexachlorobutadiene	0.238	0.220	0.214	0.204	0.212	0.218	5.9
4-Chloro-3-Methylphenol	* 0.364	0.356	0.361	0.357	0.374	0.362	2.0*
2-Methylnaphthalene	* 0.718	0.673	0.652	0.669	0.738	0.690	5.3*
Hexachlorocyclopentadiene	0.492	0.466	0.444	0.460	0.464	0.465	3.7
2,4,6-Trichlorophenol	* 0.424	0.389	0.388	0.393	0.401	0.399	3.7*
2,4,5-Trichlorophenol	* 0.444	0.418	0.382	0.372	0.345	0.392	10.0*
2-Chloronaphthalene	* 1.282	1.140	1.102	1.112	1.177	1.163	6.3*
2-Nitroaniline	0.493	0.479	0.526	0.533	0.530	0.512	4.8
Dimethyl Phthalate	1.492	1.335	1.315	1.305	1.328	1.355	5.7
Acenaphthylene	* 1.992	1.823	1.718	1.763	1.854	1.830	5.7*
2,6-Dinitrotoluene	* 0.281	0.285	0.276	0.285	0.276	0.281	1.6*
3-Nitroaniline	0.322	0.306	0.337	0.337	0.325	0.325	3.9
Acenaphthene	* 1.188	1.026	0.972	1.038	1.117	1.068	7.9*
2,4-Dinitrophenol	0.128	0.141	0.143	0.158	0.172	0.148	11.4
4-Nitrophenol	0.310	0.357	0.404	0.441	0.432	0.389	14.1
Dibenzofuran	* 1.759	1.567	1.483	1.557	1.638	1.601	6.5*
2,4-Dinitrotoluene	* 0.389	0.346	0.358	0.386	0.412	0.378	7.0*

* Compounds with required minimum RRF and maximum %RSD values.
All other compounds must meet a minimum RRF of 0.010.

SEMOVOLATILE ORGANICS INITIAL CALIBRATION DATA

ORIGINAL
(Red)Lab Name: COMPUCHEM, RTPContract: 68D10083Lab Code: COMPUCase No.: 18347

SAS No.: _____

SDG No.: CKY28Instrument ID: OWA05Calibration Date(s): 06/11/92 06/12/92Calibration Times: 2352 0249

LAB FILE ID:	RRF20 = HI920612C05	RRF50 = HG920611B05
RRF80 = HK920612C05	RRF120= HJ920612C05	RRF160= HH920612C05

COMPOUND	RRF20	RRF50	RRF80	RRF120	RRF160	RRF	% RSD
Diethylphthalate	1.657	1.593	1.522	1.560	1.578	1.582	3.1
4-Chlorophenyl-phenylether*	0.740	0.647	0.595	0.588	0.608	0.636	9.9*
Fluorene	* 1.409	1.278	1.176	1.242	1.315	1.284	6.8*
4-Nitroaniline	0.282	0.289	0.308	0.314	0.308	0.300	4.6
4,6-Dinitro-2-Methylphenol	0.133	0.147	0.147	0.142	0.137	0.141	4.4
N-Nitrosodiphenylamine (1)	0.688	0.595	0.583	0.582	0.591	0.608	7.4
4-Bromophenyl-phenylether*	0.271	0.238	0.224	0.217	0.215	0.233	9.9*
Hexachlorobenzene	* 0.296	0.259	0.243	0.242	0.239	0.256	9.3*
Pentachlorophenol	* 0.166	0.153	0.152	0.157	0.145	0.155	5.0*
Phenanthrene	* 1.231	1.091	1.033	1.105	1.140	1.120	6.5*
Anthracene	* 1.260	1.073	1.047	0.969	0.954	1.061	11.5*
Carbazole	0.994	0.887	0.857	0.850	0.786	0.875	8.7
Di-n-Butylphthalate	1.642	1.532	1.544	1.530	1.461	1.542	4.2
Fluoranthene	* 1.156	0.984	0.938	0.937	0.908	0.985	10.1*
Pyrene	* 1.750	1.584	1.479	1.324	1.439	1.515	10.6*
Ethylbenzylphthalate	0.613	0.685	0.763	0.769	0.809	0.728	10.8
3,3'-Dichlorobenzidine	0.221	0.238	0.249	0.233	0.245	0.237	4.6
Benzo(a)Anthracene	* 1.151	1.134	1.138	1.104	1.082	1.122	2.5*
Chrysene	* 1.177	1.150	1.067	0.957	0.980	1.066	9.2*
bis(2-Ethylhexyl)Phthalate	0.836	0.949	1.005	1.044	0.809	0.929	11.1
Di-n-Octyl Phthalate	1.438	1.543	1.682	1.759	1.759	1.636	8.7
Benzo(b)Fluoranthene	* 1.359	1.480	1.329	1.257	1.120	1.309	10.2*
Benzo(k)Fluoranthene	* 1.124	1.081	0.881	0.829	0.845	0.952	14.6*
Benzo(a)Pyrene	* 1.083	1.122	1.046	1.025	0.989	1.053	4.9*
Indeno(1,2,3-cd)Pyrene	* 0.626	1.081	0.992	0.918	0.897	0.903	18.9*
Dibenz(a,h)Anthracene	* 0.794	0.891	0.830	0.795	0.808	0.824	4.9*
Benzo(g,h,i)Perylene	* 0.847	0.968	0.884	0.831	0.880	0.882	6.0*
Nitrobenzene-d5	* 0.462	0.471	0.485	0.477	0.491	0.477	2.4*
2-Fluorobiphenyl	* 1.469	1.294	1.237	1.271	1.330	1.320	6.8*
Terphenyl-d14	* 1.163	1.027	0.891	0.849	0.899	0.966	13.3*
Phenol-d5	* 1.782	1.816	1.876	1.819	1.775	1.814	2.2*
2-Fluorophenol	* 1.427	1.513	1.525	1.541	1.572	1.516	3.6*
2,4,6-Tribromophenol	0.177	0.173	0.164	0.172	0.164	0.170	3.4
2-Chlorophenol-d4	* 1.327	1.295	1.315	1.294	1.292	1.305	1.2*
1,2-Dichlorobenzene-d4	* 0.969	0.916	0.907	0.911	0.936	0.928	2.8*

(1) Cannot be separated from Diphenylamine

* Compounds with required minimum RRF and maximum %RSD values.

All other compounds must meet a minimum RRF of 0.010.

7B
SEMIVOLATILE CONTINUING CALIBRATION CHECK

ORIGINAL
(Red)

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU

Case No.: 18347

SAS No.: _____

SDG No.: CKY28

Instrument ID: OWA05

Calibration date: 07/14/92 Time: 2324

Lab File ID: HG920714B05

Init. Calib. Date(s): 06/11/92 06/12/92

Init. Calib. Times: 2352 0249

COMPOUND	RRF	RRF50	MIN RRF	%D	MAX %D	Sample
Phenol	1.815	1.453	0.800	19.9	25.0	CKY29
bis(2-Chloroethyl)Ether	1.582	1.344	0.700	15.0	25.0	
2-Chlorophenol	1.463	1.289	0.800	11.9	25.0	
1,3-Dichlorobenzene	1.512	1.472	0.600	2.6	25.0	
1,4-Dichlorobenzene	1.561	1.460	0.500	6.5	25.0	
1,2-Dichlorobenzene	1.400	1.303	0.400	6.9	25.0	
2-Methylphenol	1.381	1.056	0.700	23.5	25.0	✓
2,2'-Oxybis(1-Chloropropane)	2.116	1.273		39.8		
4-Methylphenol	1.399	1.110	0.600	20.7	25.0	
N-Nitroso-Di-n-Propylamine	1.129	1.019	0.500	9.7	25.0	
Hexachloroethane	0.914	0.711	0.300	22.2	25.0	
Nitrobenzene	0.467	0.406	0.200	13.1	25.0	
Isophorone	0.877	0.758	0.400	13.6	25.0	
2-Nitrophenol	0.195	0.189	0.100	3.1	25.0	
2,4-Dimethylphenol	0.397	0.297	0.200	25.2	25.0	✓
bis(2-Chloroethoxy)Methane	0.538	0.449	0.300	16.5	25.0	
2,4-Dichlorophenol	0.308	0.308	0.200	0.0	25.0	
1,2,4-Trichlorobenzene	0.323	0.331	0.200	-2.5	25.0	
Naphthalene	1.037	0.928	0.700	10.5	25.0	
4-Chloroaniline	0.418	0.368		12.0		
Hexachlorobutadiene	0.218	0.219		-0.5		
4-Chloro-3-Methylphenol	0.362	0.335	0.200	7.5	25.0	
2-Methylnaphthalene	0.690	0.614	0.400	11.0	25.0	
Hexachlorocyclopentadiene	0.465	0.309		33.6		
2,4,6-Trichlorophenol	0.399	0.416	0.200	-4.3	25.0	
2,4,5-Trichlorophenol	0.392	0.444	0.200	-13.3	25.0	
2-Chloronaphthalene	1.163	1.109	0.800	4.6	25.0	
2-Nitroaniline	0.512	0.438		14.4		
Dimethyl Phthalate	1.355	1.356		-0.1		
Acenaphthylene	1.830	1.718	1.300	6.1	25.0	
2,6-Dinitrotoluene	0.281	0.318	0.200	-13.2	25.0	
3-Nitroaniline	0.325	0.291		10.5		
Acenaphthene	1.068	1.015	0.800	5.0	25.0	
2,4-Dinitrophenol	0.148	0.102		31.1		
4-Nitrophenol	0.389	0.147		62.2		
Dibenzofuran	1.601	1.491	0.800	6.9	25.0	
2,4-Dinitrotoluene	0.378	0.428	0.200	-13.2	25.0	

All other compounds must meet a minimum RRF of 0.010.

7C
SEMI-VOLATILE CONTINUING CALIBRATION CHECK

ORIGINAL
(Red)

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 18347 SAS No.: _____ SDG No.: CKY28

Instrument ID: OWA05 Calibration date: 07/14/92 Time: 2324

Lab File ID: HG920714B05 Init. Calib. Date(s): 06/11/92 06/12/92

Init. Calib. Times: 2352 0249

COMPOUND	RRF	RRF50	MIN RRF	%D	MAX %D
Diethylphthalate	1.582	1.447		8.5	
4-Chlorophenyl-phenylether	0.636	0.626	0.400	1.6	25.0
Fluorene	1.284	1.126	0.900	12.3	25.0
4-Nitroaniline	0.300	0.254		15.3	
4,6-Dinitro-2-Methylphenol	0.141	0.095		32.6	
N-Nitrosodiphenylamine (1)	0.608	0.499		17.9	
4-Bromophenyl-phenylether	0.233	0.230	0.100	1.3	25.0
Hexachlorobenzene	0.256	0.312	0.100	-21.9	25.0
Pentachlorophenol	0.155	0.150	0.050	3.2	25.0
Phenanthrene	1.120	0.954	0.700	14.8	25.0
Anthracene	1.061	0.944	0.700	11.0	25.0
Carbazole	0.875	0.746		14.7	
Di-n-Butylphthalate	1.542	1.343		12.9	
Fluoranthene	0.985	0.888	0.600	9.8	25.0
Pyrene	1.515	1.536	0.600	-1.4	25.0
Butylbenzylphthalate	0.728	0.775		-6.5	
3,3'-Dichlorobenzidine	0.237	0.343		44.7	
Benzo(a)Anthracene	1.122	1.102	0.800	1.8	25.0
Chrysene	1.066	1.017	0.700	4.6	25.0
bis(2-Ethylhexyl)Phthalate	0.929	1.060		-14.1	
Di-n-Octyl Phthalate	1.636	1.829		-11.8	
Benzo(b)Fluoranthene	1.309	1.066	0.700	18.6	25.0
Benzo(k)Fluoranthene	0.952	1.070	0.700	-12.4	25.0
Benzo(a)Pyrene	1.053	1.014	0.700	3.7	25.0
Indeno(1,2,3-cd)Pyrene	0.903	0.749	0.500	17.0	25.0
Dibenz(a,h)Anthracene	0.824	0.704	0.400	14.6	25.0
Benzo(g,h,i)Perylene	0.882	0.570	0.500	35.4	25.0
Nitrobenzene-d5	0.477	0.419	0.200	12.2	25.0
2-Fluorobiphenyl	1.320	1.209	0.700	8.4	25.0
Terphenyl-d14	0.966	1.016	0.500	-5.2	25.0
Phenol-d5	1.814	1.527	0.800	15.8	25.0
2-Fluorophenol	1.516	1.147	0.600	24.3	25.0
2,4,6-Tribromophenol	0.170	0.236		-38.8	
2-Chlorophenol-d4	1.305	1.177	0.800	9.8	25.0
1,2-Dichlorobenzene-d4	0.928	0.904	0.400	2.6	25.0

(1) Cannot be separated from Diphenylamine

All other compounds must meet a minimum RRF of 0.010.

6B
SEMI-VOLATILE ORGANICS INITIAL CALIBRATION DATALab Name: COMPUCHEM, RTPContract: 68D10083ORIGINAL
(Red)Lab Code: COMPUCase No.: 13347

SAS No.: _____

SDG No.: CKV14Instrument ID: OWA07Calibration Date(s): 06/15/92 06/15/92Calibration Times: 18262119

LAB FILE ID:	RRF20 = <u>HI920615B07</u>	RRF50 = <u>HG920615B07</u>
RRF80 = <u>HK920615B07</u>	RRF120= <u>HJ920615B07</u>	RRF160= <u>HH920615B07</u>

COMPOUND	RRF20	RRF50	RRF80	RRF120	RRF160	RRF	% RSD
Phenol	* 2.070	1.797	1.751	1.754	1.717	1.818	7.9*
bis(2-Chloroethyl)Ether	* 1.774	1.498	1.473	1.521	1.520	1.557	7.9*
2-Chlorophenol	* 1.759	1.454	1.426	1.406	1.341	1.477	11.0*
1,3-Dichlorobenzene	* 1.956	1.616	1.532	1.521	1.505	1.626	11.6*
1,4-Dichlorobenzene	* 1.966	1.591	1.518	1.528	1.459	1.612	12.6*
1,2-Dichlorobenzene	* 1.865	1.407	1.357	1.335	1.259	1.445	16.7*
2-Methylphenol	* 1.464	1.204	1.212	1.187	1.209	1.255	9.3*
2,2'-Oxybis(1-Chloropropane)	2.004	1.726	1.605	1.614	1.733	1.737	9.3
4-Methylphenol	* 1.516	1.350	1.333	1.340	1.332	1.374	5.8*
N-Nitroso-Di-n-Propylamine	* 1.087	0.991	0.928	0.965	0.990	0.992	5.9*
Hexachloroethane	* 0.862	0.710	0.703	0.705	0.738	0.744	9.1*
Nitrobenzene	* 0.395	0.332	0.335	0.321	0.365	0.350	8.6*
Isophorone	* 0.805	0.706	0.726	0.677	0.735	0.730	6.5*
2-Nitrophenol	* 0.235	0.198	0.200	0.195	0.202	0.206	8.0*
2,4-Dimethylphenol	* 0.392	0.312	0.307	0.294	0.331	0.327	11.8*
(2-Chloroethoxy)Methane	* 0.544	0.477	0.492	0.473	0.513	0.500	5.9*
2,4-Dichlorophenol	* 0.349	0.291	0.296	0.284	0.286	0.302	8.9*
1,2,4-Trichlorobenzene	* 0.391	0.296	0.298	0.284	0.299	0.314	13.9*
Naphthalene	* 1.309	1.006	1.000	0.882	0.919	1.024	16.4*
4-Chloroaniline	0.268	0.146	0.123	0.140	0.171	0.170	33.9
Hexachlorobutadiene	0.190	0.152	0.146	0.142	0.152	0.156	12.3
4-Chloro-3-Methylphenol	* 0.298	0.266	0.275	0.273	0.310	0.284	6.6*
2-Methylnaphthalene	* 0.849	0.701	0.716	0.665	0.665	0.719	10.6*
Hexachlorocyclopentadiene	0.346	0.334	0.315	0.327	0.356	0.336	4.8
2,4,6-Trichlorophenol	* 0.487	0.389	0.372	0.373	0.380	0.400	12.2*
2,4,5-Trichlorophenol	* 0.482	0.384	0.380	0.379	0.375	0.400	11.5*
2-Chloronaphthalene	* 1.646	1.242	1.245	1.244	1.187	1.313	14.3*
2-Nitroaniline	0.343	0.308	0.301	0.289	0.292	0.307	7.1
Dimethyl Phthalate	1.698	1.341	1.331	1.318	1.282	1.394	12.3
Acenaphthylene	* 2.522	1.943	1.951	1.820	1.810	2.009	14.6*
2,6-Dinitrotoluene	* 0.337	0.289	0.290	0.286	0.282	0.297	7.6*
3-Nitroaniline	0.111	0.048	0.020	0.014	0.028	0.044	89.7
Acenaphthene	* 1.414	1.139	1.148	1.161	1.131	1.199	10.1*
2,4-Dinitrophenol	0.115	0.120	0.126	0.138	0.146	0.129	9.9
4-Nitrophenol	0.151	0.122	0.118	0.124	0.131	0.129	10.1
Dibenzofuran	* 2.068	1.570	1.567	1.508	1.448	1.632	15.2*
2,4-Dinitrotoluene	* 0.429	0.374	0.376	0.366	0.390	0.391	5.7*

* Compounds with required minimum RRF and maximum %RSD values.
 All other compounds must meet a minimum RRF of 0.010.

6C
SEMIVOLATILE ORGANICS INITIAL CALIBRATION DATA

Original
Copy

Lab Name: COMPUCHEM, RTP Contract: 68D10083
 Lab Code: COMPU Case No.: 18347 SAS No.: SDG No.: CKY14
 Instrument ID: OWA07 Calibration Date(s): 06/15/92 06/15/92
 Calibration Times: 1826 2119

COMPOUND	RRF20	RRF50	RRF80	RRF120	RRF160	RRF	% RSD
Diethylphthalate	1.715	1.374	1.362	1.364	1.371	1.437	10.8
4-Chlorophenyl-phenylether*	0.714	0.594	0.579	0.576	0.580	0.609	9.7*
Fluorene	* 1.582	1.272	1.281	1.248	1.216	1.320	11.3*
4-Nitroaniline	0.252	0.141	0.124	0.095	0.093	0.141	46.3
4,6-Dinitro-2-Methylphenol	0.138	0.118	0.126	0.124	0.124	0.126	5.8
N-Nitrosodiphenylamine (1)	0.743	0.441	0.493	0.452	0.419	0.510	26.1
4-Bromophenyl-phenylether*	0.260	0.204	0.214	0.203	0.206	0.218	10.8*
Hexachlorobenzene	* 0.325	0.264	0.257	0.253	0.251	0.270	11.5*
Pentachlorophenol	* 0.121	0.112	0.114	0.113	0.122	0.116	4.1*
Phenanthrene	* 1.423	1.155	1.212	1.144	1.123	1.211	10.1*
Anthracene	* 1.417	1.079	1.119	1.035	0.949	1.120	15.9*
Carbazole	0.957	0.695	0.633	0.576	0.542	0.681	24.2
Di-n-Butylphthalate	1.944	1.632	1.676	1.572	1.494	1.664	10.3
Fluoranthene	* 1.195	0.986	1.027	0.965	0.956	1.030	9.3*
Pyrene	* 2.716	2.096	2.104	1.974	1.989	2.176	14.2*
Butylbenzylphthalate	1.347	1.184	1.188	1.127	1.190	1.207	6.8
3,3'-Dichlorobenzidine	0.292	0.122	0.092	0.054	0.057	0.123	79.9
Benzo(a)Anthracene	* 1.616	1.308	1.274	1.190	1.307	1.339	12.1*
Chrysene	* 1.440	1.033	1.143	1.010	1.033	1.132	15.9*
bis(2-Ethylhexyl)Phthalate	1.935	1.690	1.771	1.667	1.708	1.758	5.9
Di-n-Octyl Phthalate	2.996	2.650	2.969	2.653	2.584	2.770	7.1
Benzo(b)Fluoranthene	* 1.654	1.287	1.351	1.320	1.341	1.391	10.7*
Benzo(k)Fluoranthene	* 1.365	1.142	1.142	0.988	1.024	1.132	13.0*
Benzo(a)Pyrene	* 1.281	1.080	1.120	1.054	1.111	1.129	7.9*
Indeno(1,2,3-cd)Pyrene	* 1.434	1.088	1.170	1.123	1.069	1.177	12.6*
Dibenz(a,h)Anthracene	* 1.152	0.941	0.965	0.965	0.974	0.999	8.6*
Benzo(g,h,i)Perylene	* 1.132	0.942	0.974	0.926	0.944	0.984	8.6*
Nitrobenzene-d5	* 0.493	0.343	0.406	0.394	0.373	0.402	14.0*
2-Fluorobiphenyl	* 1.881	1.405	1.386	1.328	1.267	1.453	16.9*
Terphenyl-d14	* 1.445	1.228	1.242	1.155	1.198	1.254	8.9*
Phenol-d5	* 2.049	1.843	1.815	1.836	1.797	1.868	5.5*
2-Fluorophenol	* 1.804	1.526	1.526	1.502	1.498	1.571	8.3*
2,4,6-Tribromophenol	0.214	0.192	0.172	0.176	0.188	0.188	8.8
2-Chlorophenol-d4	* 1.583	1.314	1.290	1.273	1.233	1.339	10.4*
1,2-Dichlorobenzene-d4	* 1.218	0.938	0.924	0.905	0.848	0.967	15.0*

(1) Cannot be separated from Diphenylamine

* Compounds with required minimum RRF and maximum %RSD values.

All other compounds must meet a minimum RRF of 0.010.

73
SEMOVOLATILE CONTINUING CALIBRATION CHECK

ORIGINAL
(Red)

Lab Name: COMPUCHEM, RTP Contract: 68D10083
 Lab Code: COMPU Case No.: 18347 SAS No.: _____ SDG No.: CKY14
 Instrument ID: QWA07 Calibration date: 06/29/92 Time: 1432
 Lab File ID: HG920629A07 Init. Calib. Date(s): 06/15/92 06/15/92
 Init. Calib. Times: 1826 2119

COMPOUND	RRF	RRF50	MIN RRF	%D	MAX %D	Sample
Phenol	1.818	1.968	0.800	-8.3	25.0	CKY24
bis(2-Chloroethyl)Ether	1.557	1.568	0.700	-0.7	25.0	CKY25
2-Chlorophenol	1.477	1.437	0.800	2.7	25.0	CKY26
1,3-Dichlorobenzene	1.626	1.536	0.600	5.5	25.0	
1,4-Dichlorobenzene	1.612	1.522	0.500	5.6	25.0	
1,2-Dichlorobenzene	1.445	1.333	0.400	7.8	25.0	
2-Methylphenol	1.255	1.408	0.700	-12.2	25.0	
2,2'-Oxybis(1-Chloropropane)	1.737	1.356		21.9		
4-Methylphenol	1.374	1.455	0.600	-5.9	25.0	
N-Nitroso-Di-n-Propylamine	0.992	0.904	0.500	8.9	25.0	
Hexachloroethane	0.744	0.736	0.300	1.1	25.0	
Nitrobenzene	0.350	0.346	0.200	1.1	25.0	
Isophorone	0.730	0.721	0.400	1.2	25.0	
2-Nitrophenol	0.206	0.213	0.100	-3.4	25.0	
2,4-Dimethylphenol	0.327	0.331	0.200	-1.2	25.0	
bis(2-Chloroethoxy)Methane	0.500	0.503	0.300	-0.6	25.0	
2,4-Dichlorophenol	0.302	0.320	0.200	-6.0	25.0	
1,2,4-Trichlorobenzene	0.314	0.306	0.200	2.5	25.0	
Naphthalene	1.024	0.997	0.700	2.6	25.0	
4-Chloroaniline	0.170	0.240		-41.2		
Hexachlorobutadiene	0.156	0.170		-9.0		
4-Chloro-3-Methylphenol	0.284	0.345	0.200	-21.5	25.0	
2-Methylnaphthalene	0.719	0.695	0.400	3.3	25.0	
Hexachlorocyclopentadiene	0.336	0.289		14.0		
2,4,6-Trichlorophenol	0.400	0.410	0.200	-2.5	25.0	
2,4,5-Trichlorophenol	0.400	0.420	0.200	-5.0	25.0	
2-Chloronaphthalene	1.313	1.180	0.800	10.1	25.0	
2-Nitroaniline	0.307	0.287		6.5		
Dimethyl Phthalate	1.394	1.330		4.6		
Acenaphthylene	2.009	1.831	1.300	8.9	25.0	
2,6-Dinitrotoluene	0.297	0.307	0.200	-3.4	25.0	
3-Nitroaniline	0.044	0.199		-352.3		
Acenaphthene	1.199	1.034	0.800	13.8	25.0	
2,4-Dinitrophenol	0.129	0.121		6.2		
4-Nitrophenol	0.129	0.149		-15.5		
Dibenzofuran	1.632	1.545	0.800	5.3	25.0	
2,4-Dinitrotoluene	0.391	0.387	0.200	1.0	25.0	

All other compounds must meet a minimum RRF of 0.010.

7C
SEMIVOLATILE CONTINUING CALIBRATION CHECK

ORIGINAL
(RED)

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU

Case No.: 18347

SAS No.: _____

SDG No.: CKY14

Instrument ID: OWA07

Calibration date: 06/29/92 Time: 1432

Lab File ID: HG920629A07

Init. Calib. Date(s): 06/15/92 06/15/92

Init. Calib. Times: 1826 2119

COMPOUND	RRF	RRF50	MIN RRF	%D	MAX %D
Diethylphthalate	1.437	1.389		3.3	
4-Chlorophenyl-phenylether	0.609	0.604	0.400	0.8	25.0
Fluorene	1.320	1.130	0.900	14.4	25.0
4-Nitroaniline	0.141	0.206		-46.1	
4,6-Dinitro-2-Methylphenol	0.126	0.132		-4.8	
N-Nitrosodiphenylamine (1)	0.510	0.535		-4.9	
4-Bromophenyl-phenylether	0.218	0.246	0.100	-12.8	25.0
Hexachlorobenzene	0.270	0.284	0.100	-5.2	25.0
Pentachlorophenol	0.116	0.114	0.050	1.7	25.0
Phenanthrene	1.211	1.035	0.700	14.5	25.0
Anthracene	1.120	1.043	0.700	6.9	25.0
Carbazole	0.681	0.760		-11.6	
Di-n-Butylphthalate	1.664	1.545		7.2	
Fluoranthene	1.030	1.043	0.600	-1.3	25.0
Pyrene	2.176	1.907	0.600	12.4	25.0
Butylbenzylphthalate	1.207	1.008		16.5	
3,3'-Dichlorobenzidine	0.123	0.245		-99.2	
Benzo(a)Anthracene	1.339	1.237	0.800	7.6	25.0
Chrysene	1.132	1.103	0.700	2.6	25.0
bis(2-Ethylhexyl)Phthalate	1.758	1.438		18.2	
Di-n-Octyl Phthalate	2.770	2.290		17.3	
Benzo(b)Fluoranthene	1.391	1.352	0.700	2.8	25.0
Benzo(k)Fluoranthene	1.132	0.987	0.700	12.8	25.0
Benzo(a)Pyrene	1.129	1.079	0.700	4.4	25.0
Indeno(1,2,3-cd)Pyrene	1.177	1.051	0.500	10.7	25.0
Dibenz(a,h)Anthracene	0.999	0.919	0.400	8.0	25.0
Benzo(g,h,i)Perylene	0.984	0.879	0.500	10.7	25.0
Nitrobenzene-d5	0.402	0.366	0.200	9.0	25.0
2-Fluorobiphenyl	1.453	1.286	0.700	11.5	25.0
Terphenyl-d14	1.254	1.207	0.500	3.7	25.0
Phenol-d5	1.868	1.878	0.800	-0.5	25.0
2-Fluorophenol	1.571	1.322	0.600	15.8	25.0
2,4,6-Tribromophenol	0.188	0.175		6.9	
2-Chlorophenol-d4	1.339	1.302	0.800	2.8	25.0
1,2-Dichlorobenzene-d4	0.967	0.870	0.400	10.0	25.0

(1) Cannot be separated from Diphenylamine
All other compounds must meet a minimum RRF of 0.010.

6B
SEMIVOLATILE ORGANICS INITIAL CALIBRATION DATA

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 18347

SAS No.: _____ SDG No.: CKY14

Instrument ID: OWA02

Calibration Date(s): 06/23/92 06/23/92

Calibration Times: 1925 2154

LAB FILE ID:	RRF20 = <u>HI920623B02</u>	RRF50 = <u>HG920623B02</u>
RRF80 = <u>HK920623B02</u>	RRF120= <u>HJ920623B02</u>	RRF160= <u>HH920623B02</u>

COMPOUND	RRF20	RRF50	RRF80	RRF120	RRF160	RRF	% RSD
Phenol	* 1.864	1.592	1.623	1.450	1.493	1.604	10.1*
bis(2-Chloroethyl)Ether	* 1.695	1.490	1.488	1.354	1.392	1.484	8.9*
2-Chlorophenol	* 1.541	1.368	1.359	1.241	1.301	1.362	8.2*
1,3-Dichlorobenzene	* 1.649	1.490	1.475	1.342	1.402	1.472	7.9*
1,4-Dichlorobenzene	* 1.762	1.539	1.519	1.378	1.421	1.524	9.8*
1,2-Dichlorobenzene	* 1.567	1.398	1.395	1.263	1.287	1.382	8.7*
2-Methylphenol	* 1.296	1.122	1.122	1.018	1.050	1.122	9.6*
2,2'-Oxybis(1-Chloropropane)	1.939	1.741	1.720	1.577	1.663	1.728	7.8
4-Methylphenol	* 1.370	1.182	1.165	1.045	1.094	1.171	10.6*
N-Nitroso-Di-n-Propylamine	* 1.244	1.067	1.065	0.982	1.013	1.074	9.5*
Hexachloroethane	* 0.833	0.777	0.759	0.712	0.751	0.766	5.8*
Nitrobenzene	* 0.495	0.445	0.446	0.400	0.421	0.441	8.0*
Iscophorone	* 0.917	0.827	0.838	0.763	0.799	0.829	6.9*
2-Nitrophenol	* 0.213	0.200	0.207	0.190	0.203	0.203	4.2*
4-Dimethylphenol	* 0.382	0.327	0.345	0.303	0.336	0.339	8.5*
bis(2-Chloroethoxy)Methane	* 0.605	0.539	0.546	0.494	0.514	0.540	7.8*
2,4-Dichlorophenol	* 0.352	0.315	0.319	0.290	0.303	0.316	7.3*
1,2,4-Trichlorobenzene	* 0.384	0.349	0.339	0.307	0.323	0.340	8.6*
Naphthalene	* 1.133	1.005	0.987	0.875	0.866	0.973	11.2*
4-Chloroaniline	0.461	0.391	0.405	0.331	0.382	0.394	11.9
Hexachlorobutadiene	0.209	0.203	0.189	0.172	0.182	0.191	7.9
4-Chloro-3-Methylphenol	* 0.346	0.321	0.327	0.307	0.326	0.325	4.3*
2-Methylnaphthalene	* 0.669	0.606	0.603	0.547	0.554	0.596	8.2*
Hexachlorocyclopentadiene	0.437	0.447	0.459	0.431	0.460	0.447	2.9
2,4,6-Trichlorophenol	* 0.505	0.444	0.458	0.425	0.439	0.454	6.8*
2,4,5-Trichlorophenol	* 0.549	0.489	0.488	0.443	0.463	0.486	8.2*
2-Chloronaphthalene	* 1.475	1.288	1.313	1.193	1.203	1.294	8.8*
2-Nitroaniline	0.497	0.453	0.511	0.482	0.502	0.489	4.6
Dimethyl Phthalate	1.555	1.375	1.422	1.306	1.369	1.405	6.6
Acenaphthylene	* 2.212	1.975	1.982	1.798	1.792	1.952	8.8*
2,6-Dinitrotoluene	* 0.361	0.338	0.350	0.330	0.352	0.346	3.5*
3-Nitroaniline	0.393	0.337	0.377	0.350	0.382	0.368	6.4
Acenaphthene	* 1.332	1.174	1.177	1.069	1.089	1.168	8.9*
2,4-Dinitrophenol	0.170	0.185	0.213	0.212	0.234	0.203	12.4
4-Nitrophenol	0.159	0.150	0.169	0.159	0.173	0.162	5.6
Dibenzofuran	* 2.027	1.795	1.764	1.605	1.616	1.761	9.7*
2,4-Dinitrotoluene	* 0.480	0.464	0.485	0.454	0.477	0.472	2.7*

(1) Cannot be separated from Diphenylamine

* Compounds with required minimum RRF and maximum %RSD values.

All other compounds must meet a minimum RRF of 0.010.

6C
SEMIVOLATILE ORGANICS INITIAL CALIBRATION DATA

(Rev)

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 18347

SAS No.: _____ SDG No.: CKY14

Instrument ID: OWA02

Calibration Date(s): 06/23/92 06/23/92

Calibration Times: 1925 2154

LAB FILE ID:	RRF20 = <u>HI920623B02</u>	RRF50 = <u>HG920623B02</u>
	RRF80 = <u>HK920623B02</u>	RRF120= <u>HJ920623B02</u>
		RRF160= <u>HH920623B02</u>

COMPOUND	RRF20	RRF50	RRF80	RRF120	RRF160	RRF	% RSD
Diethylphthalate	1.578	1.457	1.528	1.432	1.480	1.495	3.9
4-Chlorophenyl-phenylether*	0.762	0.672	0.642	0.590	0.594	0.652	10.8*
Fluorene	* 1.463	1.293	1.286	1.149	1.136	1.265	10.5*
4-Nitroaniline	0.378	0.307	0.350	0.327	0.364	0.345	8.3
4,6-Dinitro-2-Methylphenol	0.153	0.161	0.179	0.175	0.189	0.171	8.4
N-Nitrosodiphenylamine (1)	0.640	0.561	0.573	0.538	0.563	0.575	6.7
4-Bromophenyl-phenylether*	0.273	0.240	0.235	0.221	0.236	0.241	8.0*
Hexachlorobenzene	* 0.360	0.317	0.308	0.295	0.308	0.318	7.9*
Pentachlorophenol	* 0.161	0.160	0.169	0.159	0.175	0.165	4.2*
Phenanthrene	* 1.296	1.141	1.164	1.073	1.101	1.155	7.5*
Anthracene	* 1.249	1.109	1.136	1.005	1.035	1.107	8.6*
Carbazole	1.013	0.892	0.964	0.897	0.928	0.939	5.4
Di-n-Butylphthalate	1.488	1.410	1.498	1.420	1.443	1.452	2.7
Fluoranthene	* 1.095	1.023	1.032	0.960	0.965	1.015	5.5*
Pyrene	* 1.897	1.645	1.744	1.582	1.723	1.718	6.9*
Butylbenzylphthalate	0.811	0.722	0.854	0.828	0.937	0.830	9.3
3,3'-Dichlorobenzidine	0.326	0.244	0.259	0.229	0.257	0.263	14.2
Benzo(a)Anthracene	* 1.393	1.171	1.162	1.101	1.147	1.195	9.5*
Chrysene	* 1.342	1.038	1.080	0.975	1.020	1.091	13.3*
bis(2-Ethylhexyl)Phthalate	1.208	1.115	1.269	1.248	1.359	1.240	7.2
Di-n-Octyl Phthalate	2.165	2.155	2.364	2.136	2.403	2.245	5.7
Benzo(b)Fluoranthene	* 1.472	1.302	1.356	1.190	1.341	1.332	7.6*
Benzo(k)Fluoranthene	* 1.353	1.194	1.140	0.989	1.196	1.174	11.1*
Benzo(a)Pyrene	* 1.224	1.030	1.040	0.898	1.071	1.053	11.1*
Indeno(1,2,3-cd)Pyrene	* 1.115	0.949	0.892	0.779	0.861	0.919	13.7*
Dibenz(a,h)Anthracene	* 0.922	0.784	0.763	0.670	0.757	0.779	11.7*
Benzo(g,h,i)Perylene	* 1.018	0.847	0.759	0.687	0.728	0.808	16.3*
Nitrobenzene-d5	* 0.485	0.443	0.456	0.411	0.428	0.445	6.3*
2-Fluorobiphenyl	* 1.612	1.379	1.373	1.230	1.245	1.368	11.2*
Terphenyl-d14	* 1.169	1.020	1.077	1.005	1.084	1.071	6.0*
Phenol-d5	* 1.821	1.614	1.615	1.471	1.523	1.609	8.3*
2-Fluorophenol	* 1.439	1.272	1.294	1.178	1.236	1.284	7.6*
2,4,6-Tribromophenol	0.239	0.229	0.229	0.212	0.233	0.228	4.4
2-Chlorophenol-d4	* 1.341	1.190	1.198	1.099	1.153	1.196	7.5*
1,2-Dichlorobenzene-d4	* 1.040	0.942	0.895	0.808	0.827	0.902	10.4*

(1) Cannot be separated from Diphenylamine

* Compounds with required minimum RRF and maximum %RSD values.

All other compounds must meet a minimum RRF of 0.010.

7B
SEMOVOLATILE CONTINUING CALIBRATION CHECK

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: CCMPU

Case No.: 18347

SAS No.: _____

SDG No.: CKY14

Instrument ID: OWA02

Calibration date: 06/29/92 Time: 1513

Lab File ID: HG920629A02

Init. Calib. Date(s): 06/23/92 06/23/92

Init. Calib. Times: 1925 2154

Sample

COMPOUND	RRF	RRF50	MIN RRF	%D	MAX %D	CKY14
Phenol	1.604	1.687	0.800	-5.2	25.0	CKY15
bis(2-Chloroethyl)Ether	1.484	1.512	0.700	-1.9	25.0	CKY16
2-Chlorophenol	1.362	1.386	0.800	-1.8	25.0	CKY17
1,3-Dichlorobenzene	1.472	1.485	0.600	-0.9	25.0	CKY18
1,4-Dichlorobenzene	1.524	1.556	0.500	-2.1	25.0	CKY18 MC
1,2-Dichlorobenzene	1.382	1.447	0.400	-4.7	25.0	CKY18 MC
2-Methylphenol	1.122	1.134	0.700	-1.1	25.0	CKY18 MC
2,2'-Oxybis(1-Chloropropane)	1.728	1.631		5.6		CKY19
4-Methylphenol	1.171	1.235	0.600	-5.5	25.0	CKY19
N-Nitroso-Di-n-Propylamine	1.074	1.267	0.500	-18.0	25.0	CKY20
Hexachloroethane	0.766	0.809	0.300	-5.6	25.0	
Nitrobenzene	0.441	0.514	0.200	-16.6	25.0	CKY22
Isophorone	0.829	0.896	0.400	-8.1	25.0	CKY23
2-Nitrophenol	0.203	0.191	0.100	5.9	25.0	
2,4-Dimethylphenol	0.339	0.338	0.200	0.3	25.0	CKY24
bis(2-Chloroethoxy)Methane	0.540	0.567	0.300	-5.0	25.0	CKY25
2,4-Dichlorophenol	0.316	0.322	0.200	-1.9	25.0	
1,2,4-Trichlorobenzene	0.340	0.355	0.200	-4.4	25.0	
Naphthalene	0.973	1.045	0.700	-7.4	25.0	
4-Chloroaniline	0.394	0.118		70.0		
Hexachlorobutadiene	0.191	0.215		-12.6		
4-Chloro-3-Methylphenol	0.325	0.336	0.200	-3.4	25.0	
2-Methylnaphthalene	0.596	0.592	0.400	0.7	25.0	
Hexachlorocyclopentadiene	0.447	0.494		-10.5		
2,4,6-Trichlorophenol	0.454	0.489	0.200	-7.7	25.0	
2,4,5-Trichlorophenol	0.486	0.508	0.200	-4.5	25.0	
2-Chloronaphthalene	1.294	1.298	0.800	-0.3	25.0	
2-Nitroaniline	0.489	0.495		-1.2		
Dimethyl Phthalate	1.405	1.447		-3.0		
Acenaphthylene	1.952	2.024	1.300	-3.7	25.0	
2,6-Dinitrotoluene	0.346	0.302	0.200	12.7	25.0	
3-Nitroaniline	0.368	0.299		18.8		
Acenaphthene	1.168	1.226	0.800	-5.0	25.0	
2,4-Dinitrophenol	0.203	0.169		16.8		
4-Nitrophenol	0.162	0.204		25.9		
Dibenzofuran	1.761	1.862	0.800	-5.7	25.0	
2,4-Dinitrotoluene	0.472	0.482	0.200	-2.1	25.0	

All other compounds must meet a minimum RRF of 0.010.

7C
SEMI VOLATILE CONTINUING CALIBRATION CHECK

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 18347 SAS No.: _____ SDG No.: CKY14

Instrument ID: OWA02 Calibration date: 06/29/92 Time: 1513

Lab File ID: HG920629A02 Init. Calib. Date(s): 06/23/92 06/23/92

Init. Calib. Times: 1925 2154

COMPOUND	RRF	RRF50	MIN RRF	%D	MAX %D
Diethylphthalate	1.495	1.512		-1.1	
4-Chlorophenyl-phenylether	0.652	0.751	0.400	-15.2	25.0
Fluorene	1.265	1.489	0.900	-17.7	25.0
4-Nitroaniline	0.345	0.289		16.2	
4,6-Dinitro-2-Methylphenol	0.171	0.157		8.2	
N-Nitrosodiphenylamine (1)	0.575	0.326		(43.3)	
4-Bromophenyl-phenylether	0.241	0.252	0.100	-4.6	25.0
Hexachlorobenzene	0.313	0.349	0.100	-9.7	25.0
Pentachlorophenol	0.165	0.186	0.050	-12.7	25.0
Phenanthrene	1.155	1.166	0.700	-1.0	25.0
Anthracene	1.107	1.055	0.700	4.7	25.0
Carbazole	0.939	0.892		5.0	
Di-n-Butylphthalate	1.452	1.498		-3.2	
Fluoranthene	1.015	1.114	0.600	-9.8	25.0
Pyrene	1.718	1.306	0.600	24.0	25.0
Butylbenzylphthalate	0.830	0.615		(25.9)	
3,3'-Dichlorobenzidine	0.263	0.165		(37.3)	
Benzo(a)Anthracene	1.195	1.190	0.800	0.4	25.0
Chrysene	1.091	0.945	0.700	13.4	25.0
bis(2-Ethylhexyl) Phthalate	1.240	1.102		11.1	
Di-n-Octyl Phthalate	2.245	2.010		10.5	
Benzo(b)Fluoranthene	1.332	1.343	0.700	-0.8	25.0
Benzo(k)Fluoranthene	1.174	1.189	0.700	-1.3	25.0
Benzo(a)Pyrene	1.053	0.979	0.700	7.0	25.0
Indeno(1,2,3-cd) Pyrene	0.919	0.862	0.500	6.2	25.0
Dibenz(a,h) Anthracene	0.779	0.729	0.400	6.4	25.0
Benzo(g,h,i) Perylene	0.808	0.785	0.500	2.8	25.0
Nitrobenzene-d5	0.445	0.525	0.200	-18.0	25.0
2-Fluorobiphenyl	1.368	1.466	0.700	-7.2	25.0
Terphenyl-d14	1.071	0.832	0.500	22.3	25.0
Phenol-d5	1.609	1.652	0.800	-2.7	25.0
2-Fluorophenol	1.284	1.406	0.600	-9.5	25.0
2,4,6-Tribromophenol	0.228	0.270		-18.4	
2-Chlorophenol-d4	1.196	1.188	0.800	0.7	25.0
1,2-Dichlorobenzene-d4	0.902	0.982	0.400	-8.9	25.0

(1) Cannot be separated from Diphenylamine
All other compounds must meet a minimum RRF of 0.010.

2C
WATER SEMIVOLATILE SURROGATE RECOVERY

OPTIONAL
RECV

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU

Case No.: 18347

SAS No.: _____

SDG No.: CKY14

EPA SAMPLE NO.	S1 (NBZ) #	S2 (FBP) #	S3 (TPH) #	S4 (PHL) #	S5 (2FP) #	S6 (TBP) #	S7 (2CP) #	S8 (DCB) #	TOT OUT
01 CKY14	92	85	133	57	88	79	109	72	0
02 CKY15	77	76	111	75	72	61	92	66	0
03 CKY16	94	88	141	92	89	74	115 *	73	1
04 CKY17	91	84	139	88	83	75	106	67	0
05 CKY18	100	90	146 *	94	89	75	115 *	75	2
06 CKY19	94	93	138	87	84	77	111 *	79	1
07 CKY20	93	92	149 *	90	86	80	109	82	1
08 CKY22	74	76	122	75	70	66	91	65	0
09 CKY23	86	85	137	83	80	76	103	70	0
10 CKY24	70	70	85	53	67	83	83	74	0
11 CKY25	62	64	82	47	56	77	72	65	0
12 CKY26	77	73	57	60	71	97	86	77	0
13 CKY18MS	71	66	108	71	66	59	85	52	0
14 CKY18MSD	0 *	0 *	4 *	2 *	2 *	2 *	2 *	0 *	8
15 SBLK13	75	68	69	67	71	84	81	68	0
16 SBLK25	74	71	108	76	70	63	92	57	0

QC LIMITS

S1 (NBZ) = Nitrobenzene-d5	(35-114)
S2 (FBP) = 2-Fluorobiphenyl	(43-116)
S3 (TPH) = Terphenyl-d14	(33-141)
S4 (PHL) = Phenol-d5	(10-110)
S5 (2FP) = 2-Fluorophenol	(21-110)
S6 (TBP) = 2,4,6-Tribromophenol	(10-123)
S7 (2CP) = 2-Chlorophenol-d4	(33-110) (advisory)
S8 (DCB) = 1,2-Dichlorobenzene-d4	(16-110) (advisory)

Column to be used to flag recovery values

* Values outside of contract required QC limits

D Surrogate diluted out

2D
SOIL SEMIVOLATILE SURROGATE RECOVERY

ORIGINAL
(Red)

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU

Case No.: 18347

SAS No.: _____

SDG No.: CKY28

Level: (low/med) LOW

EPA SAMPLE NO.	S1 (NBZ) #	S2 (FBP) #	S3 (TPH) #	S4 (PHL) #	S5 (2FP) #	S6 (TBP) #	S7 (2CP) #	S8 (DCB) #	TOT OUT
01 CKY28	73	79	79	74	79	39	92	71	0
02 CKY29	85	86	83	67	59	50	75	74	0
03 CKY30	86	76	94	80	68	65	79	83	0
04 CKY31	75	89	72	73	80	24	91	73	0
05 CKY32	81	78	98	62	39	5 * 5 *	57	78	1
06 CKY33	79	72	91	70	50	34	68	78	0
07 CKY34	25	28 *	31	25	25	21	30	26	1
08 CKY35	77	79	76	77	84	61	97	79	0
09 CKY36	72	81	80	75	79	50	96	77	0
10 CKY37	84	76	91	73	52	28	65	79	0
11 CKY38	33	38	44	33	34	32	41	31	0
12 CKY32MS	84	76	93	66	47	15 *	62	82	1
13 CKY32MSD	85	76	89	65	47	12 *	61	78	1
14 SBLK21	80	69	96	73	56	30	67	79	0
15 SBLK06	47	48	58	45	46	23	55	47	0
16 SBLK48	80	79	93	64	59	65	72	69	0

QC LIMITS

S1 (NBZ) = Nitrobenzene-d5	(23-120)
S2 (FBP) = 2-Fluorobiphenyl	(30-115)
S3 (TPH) = Terphenyl-d14	(18-137)
S4 (PHL) = Phenol-d5	(24-113)
S5 (2FP) = 2-Fluorophenol	(25-121)
S6 (TBP) = 2,4,6-Tribromophenol	(19-122)
S7 (2CP) = 2-Chlorophenol-d4	(20-130) (advisory)
S8 (DCB) = 1,2-Dichlorobenzene-d4	(20-130) (advisory)

Column to be used to flag recovery values

* Values outside of contract required QC limits

D Surrogate diluted out

2E
WATER PESTICIDE SURROGATE RECOVERY

ORIGINAL
(Rev)

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 18347 SAS No.: SDG No.: CKY14

GC Column (1): DB-608 ID: 0.53(mm) GC Column (2): RTX-1701 ID: 0.53(mm)

EPA SAMPLE NO.	TCX 1 %REC #	TCX 2 %REC #	DCB 1 %REC #	DCB 2 %REC #	OTHER (1)	OTHER (2)	TOT OUT
01 CKY14	76	87	67	53*			1
02 CKY18MS	96	86	100	95			0
03 CKY18MSD	88	86	100	87			0
04 CKY15	100	92	140	98			0
05 CKY16	83	98	76	64			0
06 CKY17	98	73	94	87			0
07 CKY18	83	84	97	81			0
08 CKY19	80	76	81	64			0
09 CKY20	84	100	85	74			0
10 CKY22	88	100	98	97			0
11 CKY23	77	83	76	81			0
12 CKY24	79	79	80	82			0
13 CKY25	86	90	91	95			0
14 CKY26	72	87	55*	80			1
15 PBLK55	99	84	130	82			0
16							
17							
18							
19							
20							
21							
22							
23							
24							
25							
26							
27							
28							
29							
30							

QC LIMITS

TCX = Tetrachloro-m-xylene (60-150)
 DCB = Decachlorobiphenyl (60-150)

Column to be used to flag recovery values.

* Values outside of QC Limits

D Surrogate diluted out.

2F
SOIL PESTICIDE SURROGATE RECOVERY

ORIGINAL
 (Red)

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 18347 SAS No.: SDG No.: CKY28

GC Column (1): DB-608 ID: 0.53(mm) GC Column (2): RTX-1701 ID: 0.53(mm)

EPA SAMPLE NO.	TCX %REC #	TCX %REC #	DCB %REC #	DCB %REC #	OTHER (1)	OTHER (2)	TOT OUT
01 CKY28	72	61	80	69			0
02 CKY29	76	64	100	88			0
03 CKY32MS	69	60	83	75			0
04 CKY32MSD	77	63	90	77			0
05 CKY30	71	53*	72	70			1
06 CKY31	80	66	100	73			0
07 CKY32	68	56*	82	69			1
08 CKY33	74	61	75	83			0
09 CKY34	75	64	88	79			0
10 CKY35	67	60	82	84			0
11 CKY36	98	59*	82	70			1
12 CKY37	71	60	72	73			0
13 CKY38	70	61	83	76			0
14 PBLK22	75	61	87	77			0
15							
16							
17							
18							
19							
20							
21							
22							
23							
24							
25							
26							
27							
28							
29							
30							

QC LIMITS

TCX = Tetrachloro-m-xylene (60-150)
 DCB = Decachlorobiphenyl (60-150)

Column to be used to flag recovery values.

* Values outside of QC Limits

D Surrogate diluted out.

WATER SEMIVOLATILE MATRIX SPIKE/MATRIX SPIKE DUPLICATE RECOVERY

Lab Name: COMPUCHEM, RTPContract: 68D10083Lab Code: COMBU Case No.: 18347

SAS No.: _____

SDG No.: CKY14ORIGINAL
(Red)Matrix Spike - EPA Sample No.: CKY18

COMPOUND	SPIKE ADDED (ug/L)	SAMPLE CONCENTRATION (ug/L)	MS CONCENTRATION (ug/L)	MS % REC #	QC LIMITS REC.
Phenol	75.00	0	49.30	66	12-110
2-Chlorophenol	75.00	0	48.80	65	27-123
1,4-Dichlorobenzene	50.00	0	25.80	52	36- 97
N-Nitroso-di-n-prop.(1)	50.00	0	29.10	58	41-116
1,2,4-Trichlorobenzene	50.00	0	23.40	57	39- 96
4-Chloro-3-methylphenol	75.00	0	52.10	69	23- 97
Acenaphthene	50.00	0	31.40	63	46-116
4-Nitrophenol	75.00	0	48.50	65	10- 80
2,4-Dinitrotoluene	50.00	0	30.80	62	24- 96
Pentachlorophenol	75.00	0	46.00	61	9-103
Pyrene	50.00	0	43.60	87	26-127

COMPOUND	SPIKE ADDED (ug/L)	MSD CONCENTRATION (ug/L)	MSD % REC #	% RPD #	QC LIMITS RPD	REC.
Phenol	75.00	1.620	2 *	188 *	42	12-110
2-Chlorophenol	75.00	1.360	2 *	188 *	40	27-123
1,4-Dichlorobenzene	50.00	0	0 *	200 *	28	36- 97
N-Nitroso-di-n-prop.(1)	50.00	0	0 *	200 *	38	41-116
1,2,4-Trichlorobenzene	50.00	0	0 *	200 *	28	39- 96
4-Chloro-3-methylphenol	75.00	1.490	2 *	189 *	42	23- 97
Acenaphthene	50.00	0	0 *	200 *	31	46-116
4-Nitrophenol	75.00	1.910	3 *	182 *	50	10- 80
2,4-Dinitrotoluene	50.00	0	0 *	200 *	38	24- 96
Pentachlorophenol	75.00	1.350	2 *	187 *	50	9-103
Pyrene	50.00	1.440	3 *	187 *	31	26-127

(1) N-Nitroso-di-n-propylamine

* Column to be used to flag recovery and RPD values with an asterisk
* Values outside of QC limitsRPD: 11 out of 11 outside limitsSpike Recovery: 11 out of 22 outside limits

COMMENTS: CLP -

CAP, HG920629A02, DI920629A02, , , ,

3D
SOIL SEMIVOLATILE MATRIX SPIKE/MATRIX SPIKE DUPLICATE RECOVERY

ORIGINAL
(Red)

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 18347

SAS No.: _____ SDG No.: CKY28

Matrix Spike - EPA Sample No.: CKY32

Level: (low/med) LOW

COMPOUND	SPIKE ADDED (ug/Kg)	SAMPLE CONCENTRATION (ug/Kg)	MS CONCENTRATION (ug/Kg)	MS % REC #	QC LIMITS REC.
Phenol	3050	0	2010	66	26- 90
2-Chlorophenol	3050	0	1661	54	25-102
4-Dichlorobenzene	2030	0	1551	76	28-104
N-Nitroso-di-n-prop.(1)	2030	0	1393	69	41-126
1,2,4-Trichlorobenzene	2030	0	1486	73	38-107
4-Chloro-3-methylphenol	3050	0	1592	52	26-103
Acenaphthene	2030	0	1531	75	31-137
4-Nitrophenol	3050	0	3411	112	11-114
2,4-Dinitrotoluene	2030	0	1665	82	28- 89
Pentachlorophenol	3050	0	1535	50	17-109
Pyrene	2030	0	1450	71	35-142

COMPOUND	SPIKE ADDED (ug/Kg)	MSD CONCENTRATION (ug/Kg)	MSD % REC #	% RPD #	QC LIMITS RPD	REC.
Phenol	3050	2026	66	0	35	26- 90
2-Chlorophenol	3050	1726	57	5	50	25-102
1,4-Dichlorobenzene	2030	1535	76	0	27	28-104
N-Nitroso-di-n-prop.(1)	2030	1466	72	4	38	41-126
1,2,4-Trichlorobenzene	2030	1499	74	1	23	38-107
4-Chloro-3-methylphenol	3050	1657	54	4	33	26-103
Acenaphthene	2030	1592	78	4	19	31-137
4-Nitrophenol	3050	3606	18 *	5	50	11-114
2,4-Dinitrotoluene	2030	1750	86	5	47	28- 89
Pentachlorophenol	3050	1466	48	4	47	17-109
Pyrene	2030	1470	72	1	36	35-142

(1) N-Nitroso-di-n-propylamine

Column to be used to flag recovery and RPD values with an asterisk

* Values outside of QC limits

RPD: 0 out of 11 outside limits

Spike Recovery: 1 out of 22 outside limits

COMMENTS: CLP
CAP, HG920701B15, DH920701B15, , , ,

EPA

United States Environmental Protection Agency
 Contract Laboratory Program Sample Management Office
 PO Box 818 Alexandria, VA 22313
 703-557-2490 FTS 557-2490

**Organic Traffic Report
 & Chain of Custody Record**
 (For Organic CLP Analysis)

SAS No.
(if applicable)

Case No.

18347

Sample Description Enter Column A)	2. Preservative (Enter In Column D)	3. Region No. Sampling Co.	5. Date Shipped	Carrier	7. Date Received -- Received by
Surface Water	III	MDE / HSWMA	6/23/92	FED EX	6/24/92 V.Balke
Ground Water		Sampler (Name)	Airbill Number		
Radial	1. HCl	JENNIFER Woods	4208457993		
Alkaline	2. HNO3				
Sol/Sediment	3. NaHSO4				
Oil/SAS	4. H2SO4				
Water (SAS)	5. Other (SAS) (Specify)				
Water (SAS)	6. Not preserved				

1. HCl	2. HNO3	3. NaHSO4	4. H2SO4	5. Other (SAS) (Specify)
6. Not preserved				

4. Type of Activity Nominal Removal

Total Pre RIFS CLEM REMA

SF Nominal RD REMA

PRPS PA RA REMA

ST SS RA REMA

FED LS O&M OIL

NPLD UST

PEYANNI TRIANGLE PARK NC 27709

ATTN: NATALIE CARTER

(714) 2-918-6735

1. Sample Identifiers (in field)	A Enter #	B Conc. Low Med High	C Sample Type: Comp./ Grab	D Preser- vative from Box 6	E RAS Analysis				F Regional Specific Tracking Number or Tag Numbers	G Station Location Number	H Mo/Day/ Year/Hour Sample Collection	I Sampler Initials	J Corresp. CLP Inorg. Samp. No.	K Sam- ple Con- dition on Arr	L High Conc. Phases (Check below)		
					VOA	BNA	Peak PCB	High ARO/ TOX							5	5	5
Y 14	3	L	G	1,6	X	X	X		3-1225137 TU	P.L.Y. 1	6/11/92 8:30	MLTY 97					
Y 15	2	L	G	1,6	X	X	X		3-1225140 TU	GW 1	6/11/92 8:30	MLTY 08					
Y 16	2	L	G	1,6	X	X	X		3-1225145 TU	GW 2	6/11/92 10:45	MLTY 04					
Y 17	2	L	G	1,6	X	X	X		3-1225146 TU	GW 3	6/11/92 12:00	MLTY 10					
Y 18	2	L	G	1,6	X	X	X		3-1225148 TU	GW 4	6/11/92 16:15	MLTY 11					
Y 19	2	L	G	1,6	X	X	X		3-1225149 TU	GW 5	6/11/92 16:45	MLTY 12					
Y 20	2	L	G	1,6	X	X	X		3-1225150 TU	GW 6	6/11/92 18:45	MLTY 13					
Y 21	2	L	G	1,6	X	X	X		3-1225151 TU								
Y 22	1	L	G	1,6	X	X	X		3-1225152 TU	GW 7	6/11/92 19:45	MLTY 15					
Y 23	1	L	G	1,6	X	X	X		3-1225153 TU	GW 8	6/11/92 21:10	MLTY 16					
Y																	

Sample used for a spike and/or duplicate
Do QC on CKY 18

Additional Sampler Signatures

Chain of Custody Seal Number
499425 499439 499444

435 440
436 441
437 442

CHAIN OF CUSTODY RECORD

Relinquished by: (Signature)	Date / Time	Received by: (Signature)	Relinquished by: (Signature)	Date / Time	Received by: (Signature)
Relinquished by: (Signature)	6/24/92 16:00		Relinquished by: (Signature)		Received by: (Signature)
Relinquished by: (Signature)			Relinquished by: (Signature)		Received by: (Signature)
Relinquished by: (Signature)			Relinquished by: (Signature)		Received by: (Signature)
Received by: (Signature)	Date / Time	Received for Laboratory by: (Signature)	Date / Time	Remarks Is custody seal intact? <input checked="" type="checkbox"/> No <input type="checkbox"/> Yes CKY 18 1-UVA RECD BROKIN THUR 3-225 16 (WED) UNCLERED <input checked="" type="checkbox"/> AIRBILL # RECD 952725505	Accuplaid (Signature)



United States Environmental Protection Agency
Contract Laboratory Program Sample Management Office
PO Box 818 Alexandria, VA 22313
703-557-2400 FTS 557-2490

**Organic Traffic Report
& Chain of Custody Record**
(For Organic CLP Analysis)

SAP
(if applicable)

Case No.

18347

Sample Description (Enter in Column A)	2. Preservative (Enter In Column D)	3. Region No. Sampling Co. Sampler (Name)	5. Date Shipped Airbill Number	Carrier	7. Date Received -- Received by Laboratory Contract Number	SAP (if applicable)	Case No.					
Surface Water	1. HCl	JENNIFER WOODS	6/13/92	FED EX	6/24/92 VDS		18347					
Ground Water	2. HNO3											
Leachate	3. NaHSO4											
Rainwater	4. H2SO4											
Solid Sediment	5. Other (SAS)											
Oil (SAS)												
Water (SAS)												
Other (BAS) (Specify)	6. Ice only											
	7. Not preserved											
CLP:	A Enter # From Box 1	B Conc. Low Med High	C Sample Type: Comp/ Grab	D Preser- vative from Box 6	E RAS Analysis	F Regional Specific Tracking Number or Tag Numbers	G Station Location Number	H Mo/Day/ Year/Time Sample Collection	I Sampler Initials	J Corresp. CLP Inorg. Samp. No.	K Sam- ple Con- dition on Rec'd	L High Conc. Phases (Check below)
KY 24	1	1	6	1,6	X	3-1225186 CO	5443	6/13/92 1050	CWT	ACLY-17		
KY 25	1	1	6	1,6	X	3-1225188 CO	5444	6/13/92 1215	JKM	ACLY-18		
KY 26	3	1	6	1,6	X	3-1225189 CO	5441	6/13/92 915	CWT	ACLY-19		
KY 27	3	1	6	1,6	X	3-1225190 CO	5440	6/13/92 1000	CWT	ACLY-20		
KY 28	5	1	6	1	X	3-1225201 CO	5421	6/13/92 945	JKM	ACLY-21		
KY 29	5	1	6	1	X	3-1225202 CO	5422	6/13/92 930	JKM	ACLY-22		
KY 30	5	1	6	1	X	3-1225203 CO	5423	6/13/92 1040	CWT	ACLY-23		
KY 31	5	1	6	1	X	3-1225210 CO	5424	6/13/92 1115	JKM	ACLY-24		
KY 32	5	1	6	1	X	3-1225211 CO	51	6/13/92 1200	T	ACLY-25		
KY 33	5	1	6	1	X	3-1225212 CO	52	6/13/92 1000	CWT	ACLY-26		
Aliquot for Cate mplete? (Y/N)	Page 1 of 3			Sample used for a spike and/or duplicate	Additional Sampler Signature Jennifer Woods			Chain of Custody Seal Number 494445 494459 494466				
Y				De QC on CKY 32				447	463			
								449	464			
								458	465			

CHAIN OF CUSTODY RECORD

Relinquished by: (Signature)	Date / Time	Received by: (Signature)	Relinquished by: (Signature)	Date / Time	Received by: (Signature)
Relinquished by: (Signature)	6/24/92 1600				
Received by: (Signature)	Date / Time	Received for Laboratory by: (Signature)	Received by: (Signature)	Date / Time	Received by: (Signature)
		V Bath			
Received by: (Signature)	Date / Time	Received for Laboratory by: (Signature)	Date / Time	Remarks	Is custody seal intact? <input checked="" type="checkbox"/> None
			6/24/92 0830	RECV'D CKY 31 1-VAT BROKEN IN BOX TO SAUCL TAB 4 3-1225210 6/24/92 AIRBILL # 2957725505	
Split Samples <input type="checkbox"/> Accepted (Signature)					



COMPUCHEM
LABORATORIES, INC.

P.O. Box 12652 3308 Chapel Hill/Nelson Highway Research Triangle Park, NC 27709 (919) 549-8263

ORIGINAL
(Red)

SDG NARRATIVE SDG # CKY14
CASE # 18347
CONTRACT # 68D10083

SAMPLES: CKY14, CKY15, CKY16, CKY17, CKY18, CKY19, CKY20, CKY22,
CKY23, CKY24, CKY25, CKY26

This portion of Case # 18347 consisted of twelve liquid samples for volatile, semivolatile, and pesticide analysis. They were received on 6/24/92 via Federal Express in properly sealed shipping containers with traffic reports. One volatile container of CKY16 was received broken.

VOLATILES:

All volatile fractions were analyzed within holding time requirements. Samples containing no target compounds above the CRQL were CKY15, CKY17, and CKY26. The remaining samples contained methylene chloride and/or acetone, which were also detected in the associated instrument blanks. All system monitoring compounds met recovery criteria. The matrix spike/matrix spike duplicate results were acceptable. The raw sample pH values are given on the attached pages.

SEMICVOLATILES:

All semivolatile fractions were extracted and analyzed within holding time requirements. None of the samples contained any target compounds above the CRQL. The recovery of D14-terphenyl exceeded QC limits in samples CKY13 and CKY20. All other production sample surrogates met recovery criteria. The matrix spike and blank spike met overall acceptance criteria. However, the matrix spike duplicate results failed both surrogate and spike recovery criteria due to an extraction error. Since this was not discovered until holding times had expired and since the other recoveries were acceptable, no reextraction were performed.

PESTICIDES:

All pesticide fractions were extracted and analyzed within holding time requirements. None of the samples contained any target compounds above the CRQL. All surrogate recoveries were acceptable. The matrix spike/matrix spike duplicate results were acceptable. Endrin ketone, which was detected in the spikes but not the original, was identified as a spike compound degradation product. All associated sequences met degradation criteria.



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I certify that this data package is in compliance with the terms and conditions of the contract, both technically and for completeness, for other than the conditions detailed above. Release of the data contained in this hardcopy data package and in the computer-readable data submitted on floppy diskette has been authorized by the Laboratory Manager or his designee, as verified by the following signature.

Note: This report was paginated for reference and accountability in increasing numerical sequence.

Sarah A. Hubbard
Sarah A. Hubbard 7/13/92
Technical Reviewer



COMPUCHEM
LABORATORIES, INC.

P.O. Box 12652 3308 Chapel Hill / Nelson Highway Research Triangle Park, NC 27709 (919) 549-8263

ORIGINAL
(Red)

SDG NARRATIVE -- CASE 18347
SDG NO. CKY28
Contract No. 68D10083
Compuchem Laboratories, Inc.

Samples: CKY28, CKY29, CKY30, CKY31, CKY32, CKY33, CKY34, CKY35,
CKY36, CKY37, CKY38

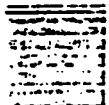
Attached are pertinent Quality Assurance Notices dealing with the analysis of eleven (11) soil samples associated with Case 18347, SDG No. CKY28. With one exception, the samples were received intact on June 24, 1992 in properly sealed shipping containers with the corresponding chain-of-custody documents and traffic reports. For sample CKY31, one volatile container was received broken. The courier was Federal Express. The samples were logged into the Compuchem Laboratory Management system and scheduled for the analysis of the volatile, semi-volatile, and pesticide/PCB fractions. The percent moisture content values of the samples ranged from eight (8.0) to forty (40.0). The pH values of the samples ranged from 5.5 to 7.2, and were within the limits of 5.0 to 11.0 as specified by EPA protocols. This SDG narrative addresses the volatile and the semi-volatile fractions.

VOLATILES

The samples were prepared and analyzed within the proper holding time requirements. EPA target compound list (TCL) analytes methylene chloride and acetone were detected in all of the samples. Toluene was also detected in sample CKY28. The concentration of toluene in sample CKY28 was below the contract required quantitation limit (CRQL). Consequently, this analyte has been flagged with the 'J' qualifier on the organic analysis data sheets (CADS). Tentatively identified compounds (TICs) were detected in all of the samples ranging in number per sample from one (1) to two (2). The TICs detected in the samples were assessed as laboratory artifacts. A laboratory artifact is usually an early eluting peak believed to consist of water and various atmospheric gases. Quality assurance notices have been included with the data for the laboratory artifacts.

In the volatile fraction, recovery and RPD values met QC limits in the matrix spike, CKY32 MS, and the matrix spike duplicate, CKY32 MSD. In addition to the spiking compounds, methylene chloride and acetone were detected in the MS and the MSD. Methylene chloride and acetone were detected in all of the four associated volatile method blanks. Each of the blanks contained one (1) TIC. The TIC found in each of the method blanks was assessed as a laboratory artifact. System monitoring compound recovery values for the samples, the blanks, and the duplicate sample spikes passed contract required QC limits.

Results of the original sample, CKY32, and the MS did not agree favorably. Variations were observed in the number of extraneous



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ORIGINAL
(Red)

peaks. These variations can be attributed to the inhomogeneity of the soil matrix. We are reporting the data for the original sample and the MS with reference to the enclosed quality assurance notice.

SEMI-VOLATILES

Initially, all of the samples were extracted and analyzed within holding time limits. In the initial extraction of samples CKY28, CKY29, CKY31, CKY34, CKY35, CKY36, and CKY38, performed within holding time limits, surrogate recovery values did not meet QC criteria. The samples were re-extracted outside of holding time requirements. Although performed outside of holding time limits, extraction efficiency was improved and overall surrogate recovery criteria were met for these samples. Results were comparable between the two analyses for each of the samples. The data for the original analyses have been included in the non-deliverable section of the complete SDG file (CSF). We are reporting the data from the second extraction of these samples with reference to the enclosed quality assurance notice.

Samples CKY28, CKY31, and CKY36 did not contain any TCL compounds. TCL analytes were detected in the remaining samples ranging in number per sample from one (1) to twelve (12). These analytes included 4-methylphenol, phenanthrene, butylbenzylphthalate, pyrene, fluoranthene, bis(2-ethylhexyl)phthalate, 2-methylnaphthalene, benzo(a)anthracene, chrysene, benzo(a)pyrene, indeno(1,2,3-cd)pyrene, and benzo(b/k)fluoranthene. Benzo(b/k) fluoranthene coeluted in samples CKY29 and CKY38. Extracted ion current profiles (EICP) have been included with samples CKY29 and CKY38 where the fluoranthene isomers coeluted. TICs were detected in all of the samples ranging in number per sample from six (6) to twenty-six (26). Some of the TICs detected in the samples were assessed as blank contaminants and laboratory artifacts. Quality assurance notices have been included with the data for the samples where the laboratory artifacts were detected.

Extensive internal tests performed here at Compuchem indicated that a particular lot of methylene chloride contained several early eluting TICs. These TICs appeared in samples even though a pre-approval test performed on the new lot did not contain the TICs. Samples CKY28, CKY29, CKY30, CKY31, CKY32, CKY33, CKY34, CKY35, CKY36, CKY37, CKY38, and method blank S6LK21 were affected by this problem with the methylene chloride. The discovery of the contaminated methylene chloride was not made until after these samples and the method blank had been extracted. We have received a new lot of solvent which is contaminant free. We are reporting the data for these samples and the method blank with reference to the enclosed laboratory notices.

In the semi-volatile fraction, with one exception, recovery and RPD values met QC limits in the matrix spike, CKY32 MS, and the matrix spike duplicate, CKY32 MSD. Spiking compound 4-nitrophenol exceeded the QC limits. In addition to the spiking compounds,



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ORIGINAL
(Red)

butylbenzylphthalate was detected in the MS and the MSD. Benzo(b/k)fluoranthene coeluted in the MSD. Method blank SBLK06 did not contain any TCL compounds. Butylbenzylphthalate was detected below the CRQL in method blanks SBLK21 and SBLK48. There were three TICs detected in method blank SBLK06. Method blank SBLK21 contained six (6) TICs. In method blank SBLK48, five (5) TICs were detected.

Surrogate 2,4,6-tribromophenol failed QC limits in sample CKY32 and the duplicate sample spikes, CKY32MS/MSD. The unacceptable recoveries in the matrix spikes and the original sample have been attributed to the particular sample matrix rather than to deficiencies in the laboratory's analytical system. We are reporting the data for the original sample and the matrix spikes with reference to the enclosed quality assurance notice. In sample CKY34, 2-fluorobiphenyl failed QC limits. The remaining surrogate recovery values for the samples, the blanks, and the sample spikes passed contract required QC limits.

I certify that this data package is in compliance with the terms and conditions of the contract, both technically and for completeness, for other than the conditions detailed above. Release of the data contained in this hardcopy data package and in the computer-readable data submitted on floppy diskette has been authorized by the Laboratory Manager or his designee, as verified by the following signature:

Cynthia E. Edwards 07/17/92
Cynthia E. McCloud- Edwards
Technical Reviewer
July 10, 1992

Note: This report is paginated for reference and accountability.



COMPUCHEM
LABORATORIES, INC.

P.O. Box 12652 3308 Chapel Hill/Nelson Highway Research Triangle Park, NC 27709 (919) 549-8263

ORIGINAL
(Red)

SDG NARRATIVE -- CASE 18347

SDG NO. CKY28

Contract No. 68D10083

CompuChem Laboratories, Inc.

Samples: CKY28, CKY29, CKY30, CKY31, CKY32, CKY33, CKY34, CKY35,
CKY36, CKY37, CKY38

Pesticides/PCBs

The samples were extracted and analyzed within the proper holding time requirements. No TCL compounds were confirmed in the samples at concentrations above the CRQL. However, TCL compounds were confirmed in the samples at concentrations below the CRQL. These compounds included alpha-BHC, gamma-BHC, heptachlor, aldrin, heptachlor epoxide, endosulfan I, endosulfan II, dieldrin, endrin, 4,4'-DDD, 4,4'-DDE, 4,4'-DDT, methoxychlor, endosulfan sulfate, alpha-chlordane, gamma-chlordane, and endrin ketone. The "p" qualifier was applied to some of the compounds confirmed in these samples. The "p" flag was used to indicate that the differences between the detected concentrations for these compounds analyzed on the DB-608 column and the RTX-1701 column were greater than twenty-five (25) percent. TCL compound methoxychlor was confirmed in the method blank at a concentration below the CRQL.

Recovery and RPD values met QC limits in the matrix spike, CKY32MS, and the matrix spike duplicate, CKY32MSD.

The recovery of the tetrachloro-m-xylene surrogate for samples CKY30, CKY32, and CKY36 analyzed on the RTX-1701 column was below advisory QC limits.

I certify that this data package is in compliance with the terms and conditions of the contract, both technically and for completeness, for other than the conditions detailed above. Release of the data contained in this hardcopy data package and in the computer readable floppy diskette has been authorized by the Laboratory Manager or his designee, as verified by the following signature:

Jean M. Zimmerman
Jean M. Zimmerman
Technical Reviewer
July 16, 1992

Note: This report is paginated for reference and accountability.

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

(Rev)

CKY18MS

Lab Name: COMPUCHEM, RTPContract: 68D10083Lab Code: COMPU Case No.: 18347SAS No.: _____ SDG No.: CKY14Matrix: (soil/water) WATERLab Sample ID: 499426Sample wt/vol: 5.0 (g/mL) MLLab File ID: CN099426B56Level: (low/med) LOWDate Received: 06/24/92

% Moisture: not dec. _____

Date Analyzed: 06/25/92Column: DB624 ID: 0.530 (mm)Dilution Factor: 1.0

Soil Extract Volume: _____ (uL)

Soil Aliquot Volume: _____ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS:	
		(ug/L or ug/Kg)	UG/L
74-87-3-----	Chloromethane	10	U
74-83-9-----	Bromomethane	10	U
75-01-4-----	Vinyl Chloride	10	U
75-00-3-----	Chloroethane	10	U
75-09-2-----	Methylene Chloride	12	B
67-64-1-----	Acetone	10	U
75-15-0-----	Carbon Disulfide	10	U
75-35-4-----	1,1-Dichloroethene	57	
75-34-3-----	1,1-Dichloroethane	10	U
540-59-0-----	1,2-Dichloroethene (total)	10	U
67-66-3-----	Chloroform	10	U
107-06-2-----	1,2-Dichloroethane	10	U
78-93-3-----	2-Butanone	10	U
71-55-6-----	1,1,1-Trichloroethane	10	U
56-23-5-----	Carbon Tetrachloride	10	U
75-27-4-----	Bromodichloromethane	10	U
78-87-5-----	1,2-Dichloroproppane	10	U
10061-01-5-----	cis-1,3-Dichloropropene	10	U
79-01-6-----	Trichloroethene	54	
124-48-1-----	Dibromoethylmethane	10	U
79-00-5-----	1,1,2-Trichloroethane	10	U
71-43-2-----	Benzene	55	
10061-02-6-----	Trans-1,3-Dichloropropene	10	U
75-25-2-----	Bromoform	10	U
108-10-1-----	4-Methyl-2-Pentanone	10	U
591-78-6-----	2-Hexanone	10	U
127-18-4-----	Tetrachloroethene	10	U
79-34-5-----	1,1,2,2-Tetrachloroethane	10	U
108-88-3-----	Toluene	53	
108-90-7-----	Chlorobenzene	53	
100-41-4-----	Ethylbenzene	10	U
100-42-5-----	Styrene	10	U
1330-20-7-----	Xylene (total)	10	U

CAS NO.	COMPOUND	CONCENTRATION UNITS:	Q
74-87-3-----	Chloromethane	10	U
74-83-9-----	Bromomethane	10	U
75-01-4-----	Vinyl Chloride	10	U
75-00-3-----	Chloroethane	10	U
75-09-2-----	Methylene Chloride	12	B
67-64-1-----	Acetone	10	U
75-15-0-----	Carbon Disulfide	10	U
75-35-4-----	1,1-Dichloroethene	57	
75-34-3-----	1,1-Dichloroethane	10	U
540-59-0-----	1,2-Dichloroethene (total)	10	U
67-66-3-----	Chloroform	10	U
107-06-2-----	1,2-Dichloroethane	10	U
78-93-3-----	2-Butanone	10	U
71-55-6-----	1,1,1-Trichloroethane	10	U
56-23-5-----	Carbon Tetrachloride	10	U
75-27-4-----	Bromodichloromethane	10	U
78-87-5-----	1,2-Dichloroproppane	10	U
10061-01-5-----	cis-1,3-Dichloropropene	10	U
79-01-6-----	Trichloroethene	54	
124-48-1-----	Dibromoethylmethane	10	U
79-00-5-----	1,1,2-Trichloroethane	10	U
71-43-2-----	Benzene	55	
10061-02-6-----	Trans-1,3-Dichloropropene	10	U
75-25-2-----	Bromoform	10	U
108-10-1-----	4-Methyl-2-Pentanone	10	U
591-78-6-----	2-Hexanone	10	U
127-18-4-----	Tetrachloroethene	10	U
79-34-5-----	1,1,2,2-Tetrachloroethane	10	U
108-88-3-----	Toluene	53	
108-90-7-----	Chlorobenzene	53	
100-41-4-----	Ethylbenzene	10	U
100-42-5-----	Styrene	10	U
1330-20-7-----	Xylene (total)	10	U

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

ORIGINAL
CKY18MSD

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 18347

SAS No.: _____

SDG No.: CKY14

Matrix: (soil/water) WATER

Lab Sample ID: 499427

Sample wt/vol: 5.0 (g/mL) ML

Lab File ID: CN099427B56

Level: (low/med) LOW

Date Received: 06/24/92

% Moisture: not dec.

Date Analyzed: 06/25/92

GC Column: DB624 ID: 0.530 (mm)

Dilution Factor: 1.0

Soil Extract Volume: _____ (uL)

Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

(ug/L or ug/Kg) UG/L

CAS NO.

COMPOUND

Q

74-87-3-----	Chloromethane	10	U
74-83-9-----	Bromomethane	10	U
75-01-4-----	Vinyl Chloride	10	U
75-00-3-----	Chloroethane	10	U
75-09-2-----	Methylene Chloride	9	BJ
67-64-1-----	Acetone	11	B
75-15-0-----	Carbon Disulfide	10	U
75-35-4-----	1,1-Dichloroethene	56	
75-34-3-----	1,1-Dichloroethane	10	U
540-59-0-----	1,2-Dichloroethene (total)	10	U
67-66-3-----	Chloroform	10	U
107-06-2-----	1,2-Dichloroethane	10	U
78-93-3-----	2-Butanone	10	U
71-55-6-----	1,1,1-Trichloroethane	10	U
56-23-5-----	Carbon Tetrachloride	10	U
75-27-4-----	Bromodichloromethane	10	U
78-87-5-----	1,2-Dichloropropane	10	U
10061-01-5-----	cis-1,3-Dichloropropene	10	U
79-01-6-----	Trichloroethene	56	
124-48-1-----	Dibromochloromethane	10	U
79-00-5-----	1,1,2-Trichloroethane	10	U
71-43-2-----	Benzene	57	
10061-02-6-----	Trans-1,3-Dichloropropene	10	U
75-25-2-----	Bromoform	10	U
108-10-1-----	4-Methyl-2-Pentanone	10	U
591-78-6-----	2-Hexanone	10	U
127-18-4-----	Tetrachloroethene	10	U
79-34-5-----	1,1,2,2-Tetrachloroethane	10	U
108-88-3-----	Toluene	55	
108-90-7-----	Chlorobenzene	55	
100-41-4-----	Ethylbenzene	10	U
100-42-5-----	Styrene	10	U
1330-20-7-----	Xylene (total)	10	U

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CKY32MS

Lab Name: COMPUCHEM, RTPContract: 68D10083Lab Code: COMPUCase No.: 18347

SAS No.: _____

SDG No.: CKY28Matrix: (soil/water) SOILLab Sample ID: 499453Sample wt/vol: 5.0 (g/mL) GLab File ID: G2R99453A54Level: (low/med) LOWDate Received: 06/24/92% Moisture: not dec. 18Date Analyzed: 06/26/92GC Column: DB-624 ID: 0.530 (mm)Dilution Factor: 1.0

Soil Extract Volume: _____ (uL)

Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

(ug/L or ug/Kg) UG/KG

Q

CAS NO.	COMPOUND	UG/KG	Q
74-87-3-----	Chloromethane	12	U
74-33-9-----	Bromomethane	12	U
75-01-4-----	Vinyl Chloride	12	U
75-00-3-----	Chloroethane	12	U
75-09-2-----	Methylene Chloride	32	B
67-64-1-----	Acetone	29	B
75-15-0-----	Carbon Disulfide	12	U
75-35-4-----	1,1-Dichloroethene	71	
75-34-3-----	1,1-Dichloroethane	12	U
540-59-0-----	1,2-Dichloroethene (total)	12	U
67-66-3-----	Chloroform	12	U
107-06-2-----	1,2-Dichloroethane	12	U
78-93-3-----	2-Butanone	12	U
71-55-6-----	1,1,1-Trichloroethane	12	U
56-23-5-----	Carbon Tetrachloride	12	U
75-27-4-----	Bromodichloromethane	12	U
78-37-5-----	1,2-Dichloropropane	12	U
10061-01-5-----	cis-1,3-Dichloropropene	12	U
79-01-6-----	Trichloroethene	64	
124-48-1-----	Dibromoethylmethane	12	U
79-00-5-----	1,1,2-Trichloroethane	12	U
71-43-2-----	Benzene	70	
10061-02-6-----	Trans-1,3-Dichloropropene	12	U
75-25-2-----	Bromoform	12	U
108-10-1-----	4-Methyl-2-Pentanone	12	U
591-78-6-----	2-Hexanone	12	U
127-18-4-----	Tetrachloroethene	12	U
79-34-5-----	1,1,2,2-Tetrachloroethane	12	U
108-88-3-----	Toluene	69	
108-90-7-----	Chlorobenzene	65	
100-41-4-----	Ethylbenzene	12	U
100-42-5-----	Styrene	12	U
1330-20-7-----	Xylene (total)	12	U

1A
VOLATILE ORGANICS ANALYSIS DATA SHEETOR EPA SAMPLE NO.
(Ref)

CKY32MSD

Lab Name: COMPUCHEM, RTPContract: 68D10083Lab Code: COMPU Case No.: 18347SAS No.: _____ SDG No.: CKY28Matrix: (soil/water) SOILLab Sample ID: 499454Sample wt/vol: 5.0 (g/mL) GLab File ID: GRO99454A54Level: (low/med) LOWDate Received: 06/24/92% Moisture: not dec. 18Date Analyzed: 06/25/92GC Column: DB-624 ID: 0.530 (mm)Dilution Factor: 1.0

Soil Extract Volume: _____ (uL)

Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

(ug/L or ug/Kg) UG/KG

Q

CAS NO.

COMPOUND

74-37-3-----	Chloromethane	12	U
74-83-9-----	Bromomethane	12	U
75-01-4-----	Vinyl Chloride	12	U
75-00-3-----	Chloroethane	12	U
75-09-2-----	Methylene Chloride	38	B
67-64-1-----	Acetone	30	B
75-15-0-----	Carbon Disulfide	12	U
75-35-4-----	1,1-Dichloroethene	82	
75-34-3-----	1,1-Dichloroethane	12	U
540-59-0-----	1,2-Dichloroethene (total)	12	U
67-66-3-----	Chloroform	12	U
107-06-2-----	1,2-Dichloroethane	12	U
78-93-3-----	2-Butanone	12	U
71-55-6-----	1,1,1-Trichloroethane	12	U
56-23-5-----	Carbon Tetrachloride	12	U
75-27-4-----	Bromodichloromethane	12	U
78-87-5-----	1,2-Dichloropropane	12	U
10061-01-5-----	cis-1,3-Dichloropropene	12	U
79-01-6-----	Trichloroethene	74	
124-48-1-----	Dibromo-chloromethane	12	U
79-00-5-----	1,1,2-Trichloroethane	12	U
71-43-2-----	Benzene	80	
10061-02-6-----	Trans-1,3-Dichloropropene	12	U
75-25-2-----	Bromoform	12	U
108-10-1-----	4-Methyl-2-Pentanone	12	U
591-78-6-----	2-Hexanone	12	U
127-18-4-----	Tetrachloroethene	12	U
79-34-5-----	1,1,2,2-Tetrachloroethane	12	U
108-88-3-----	Toluene	79	
108-90-7-----	Chlorobenzene	78	
100-41-4-----	Ethylbenzene	12	U
100-42-5-----	Styrene	12	U
1330-20-7-----	Xylene (total)	12	U

1B
SEMICVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

ORIGINAL
Recky18MSLab Name: COMPUCHEM, RTP Contract: 68D10083Lab Code: COMPU Case No.: 18347 SAS No.: _____ SDG No.: CKY14Matrix: (soil/water) WATER Lab Sample ID: 499429Sample wt/vol: 500 (g/mL) ML Lab File ID: GH099429B02Level: (low/med) LOW Date Received: 06/24/92% Moisture: _____ decanted: (Y/N) _____ Date Extracted: 06/26/92Concentrated Extract Volume: 500.0 (uL) Date Analyzed: 06/29/92Injection Volume: 2.0(uL) Dilution Factor: 1.0GPC Cleanup: (Y/N) N pH: _____CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/L

Q

CAS NO.	COMPOUND	Q
108-95-2	Phenol	49
111-44-4	bis(2-Chloroethyl)Ether	10
95-57-8	2-Chlorophenol	49
541-73-1	1,3-Dichlorobenzene	10
106-46-7	1,4-Dichlorobenzene	26
95-50-1	1,2-Dichlorobenzene	10
95-48-7	2-Methylphenol	10
108-60-1	2,2'-Oxybis(1-Chloropropane)	10
106-44-5	4-Methylphenol	10
621-64-7	N-Nitroso-Di-n-Propylamine	29
67-72-1	Hexachloroethane	10
98-95-3	Nitrobenzene	10
78-59-1	Iscophorone	10
88-75-5	2-Nitrophenol	10
105-67-9	2,4-Dimethylphenol	10
111-91-1	bis(2-Chloroethoxy)Methane	10
120-83-2	2,4-Dichlorophenol	10
120-32-1	1,2,4-Trichlorobenzene	28
91-20-3	Naphthalene	10
106-47-6	4-Chloroaniline	10
87-68-3	Hexachlorobutadiene	10
59-50-7	4-Chloro-3-Methylphenol	52
91-57-6	2-Methylnaphthalene	10
77-47-4	Hexachlorocyclopentadiene	10
88-06-2	2,4,6-Trichlorophenol	10
95-95-4	2,4,5-Trichlorophenol	25
91-58-7	2-Chloronaphthalene	10
88-74-4	2-Nitroaniline	25
131-11-3	Dimethyl Phthalate	10
208-96-8	Acenaphthylene	10
606-20-2	2,6-Dinitrotoluene	10
99-09-2	3-Nitroaniline	25
83-32-9	Acenaphthene	31

1C
SEMICVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

ORIGINAL
(Rev)

CKY18MS

Lab Name: COMPUCHEM, RTPContract: 68D10083Lab Code: COMPU Case No.: 18347SAS No.: _____ SDG No.: CKY14Matrix: (soil/water) WATERLab Sample ID: 499429Sample wt/vol: 500 (g/mL) MLLab File ID: GH099429B02Level: (low/med) LOWDate Received: 06/24/92

% Moisture: _____ decanted: (Y/N) _____

Date Extracted: 06/26/92Concentrated Extract Volume: 500.0 (uL)Date Analyzed: 06/29/92Injection Volume: 2.0 (uL)Dilution Factor: 1.0GPC Cleanup: (Y/N) N pH: _____CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/L Q

CAS NO. COMPOUND

51-28-5-----	2,4-Dinitrophenol	25	U
100-02-7-----	4-Nitrophenol	48	
132-64-9-----	Dibenzofuran	10	U
121-14-2-----	2,4-Dinitrotoluene	31	
84-66-2-----	Diethylphthalate	10	U
7005-72-3-----	4-Chlorophenyl-phenylether	10	U
86-73-7-----	Fluorene	10	U
100-01-6-----	4-Nitroaniline	25	U
534-52-1-----	4,6-Dinitro-2-Methylphenol	25	U
86-30-6-----	N-Nitrosodiphenylamine (1)	10	U
101-55-3-----	4-Bromophenyl-phenylether	10	U
118-74-1-----	Hexachlorobenzene	10	U
87-86-5-----	Pentachlorophenol	46	
85-01-8-----	Phenanthrene	10	U
120-12-7-----	Anthracene	10	U
86-74-8-----	Carbazole	10	U
84-74-2-----	Di-n-Butylphthalate	10	U
206-44-0-----	Fluoranthene	10	U
129-00-0-----	Pyrene	44	
85-68-7-----	Butylbenzylphthalate	10	U
91-94-1-----	3,3'-Dichlorobenzidine	10	U
56-55-3-----	Benz(a)Anthracene	10	U
218-01-9-----	Chrysene	10	U
117-81-7-----	bis(2-Ethylhexyl) Phthalate	2	BJ
117-84-0-----	Di-n-Octyl Phthalate	10	U
205-99-2-----	Benz(b)Fluoranthene	10	U
207-08-9-----	Benz(k)Fluoranthene	10	U
50-32-8-----	Benz(a)Pyrene	10	U
193-39-5-----	Indeno(1,2,3-cd) Pyrene	10	U
53-70-3-----	Dibenz(a,h)Anthracene	10	U
191-24-2-----	Benzo(g,h,i)Perylene	10	U

(1) - Cannot be separated from Diphenylamine

1B
SEMICVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

09/12/92
(Rev)

CKY18MSD

Lab Name: COMPUCHEM, RTPContract: 68D10083Lab Code: COMPUCase No.: 18347

SAS No.: _____

SDG No.: CKY14Matrix: (soil/water) WATERLab Sample ID: 499430Sample wt/vol: 500 (g/mL) MLLab File ID: GH099430B02Level: (low/med) LOWDate Received: 06/24/92

Moisture: _____ decanted: (Y/N) _____

Date Extracted: 06/26/92Concentrated Extract Volume: 500.0 (uL)Date Analyzed: 06/29/92Injection Volume: 2.0 (uL)Dilution Factor: 1.0GPC Cleanup: (Y/N) N pH: _____

CONCENTRATION UNITS:

(ug/L or ug/Kg) UG/LQ

CAS NO.	COMPOUND	2	BJ
108-95-2-----	Phenol	10	U
111-44-4-----	bis(2-Chloroethyl)Ether	1	J
95-57-8-----	2-Chlorophenol	10	U
541-73-1-----	1,3-Dichlorobenzene	10	U
106-46-7-----	1,4-Dichlorobenzene	10	U
95-50-1-----	1,2-Dichlorobenzene	10	U
95-48-7-----	2-Methylphenol	10	U
108-60-1-----	2,2'-Oxybis(1-Chloropropane)	10	U
106-44-5-----	4-Methylphenol	10	U
621-64-7-----	N-Nitroso-Di-n-Propylamine	10	U
67-72-1-----	Hexachloroethane	10	U
98-95-3-----	Nitrobenzene	10	U
78-59-1-----	Isophorone	10	U
88-75-5-----	2-Nitrophenol	10	U
105-67-9-----	2,4-Dimethylphenol	10	U
111-91-1-----	bis(2-Chloroethoxy)Methane	10	U
120-83-2-----	2,4-Dichlorophenol	10	U
120-82-1-----	1,2,4-Trichlorobenzene	10	U
91-20-3-----	Naphthalene	10	U
106-47-8-----	4-Chloroaniline	10	U
87-68-3-----	Hexachlorobutadiene	10	U
59-50-7-----	4-Chloro-3-Methylphenol	1	J
91-57-6-----	2-Methylnaphthalene	10	U
77-47-4-----	Hexachlorocyclopentadiene	10	U
88-06-2-----	2,4,6-Trichlorophenol	10	U
95-95-4-----	2,4,5-Trichlorophenol	25	U
91-58-7-----	2-Chloronaphthalene	10	U
88-74-4-----	2-Nitroaniline	25	U
131-11-3-----	Dimethyl Phthalate	10	U
208-96-8-----	Acenaphthylene	10	U
606-20-2-----	2,6-Dinitrotoluene	10	U
99-09-2-----	3-Nitroaniline	25	U
63-32-9-----	Acenaphthene	10	U

1C
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

ORIGINAL
(Red)

CKY18MSD

Lab Name: COMPUCHEM, RTP Contract: 68D10083Lab Code: COMPU Case No.: 18347 SAS No.: _____ SDG No.: CKY14Matrix: (soil/water) WATERLab Sample ID: 499430Sample wt/vol: 500 (g/mL) MLLab File ID: GH099430B02Level: Low (low/med) LOWDate Received: 06/24/92Moisture: _____ decanted: (Y/N) Date Extracted: 06/26/92Concentrated Extract Volume: 500.0 (uL)Date Analyzed: 06/29/92Injection Volume: 2.0 (uL)Dilution Factor: 1.0GPC Cleanup: (Y/N) N pH: CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/L Q

CAS NO.	COMPOUND	UG/L	Q
51-28-5	2,4-Dinitrophenol	25	U
100-02-7	4-Nitrophenol	2	J
132-64-9	Dibenzofuran	10	U
121-14-2	2,4-Dinitrotoluene	10	U
84-66-2	Diethylphthalate	10	U
7005-72-3	4-Chlorophenyl-phenylether	10	U
86-73-7	Fluorene	10	U
100-01-6	4-Nitroaniline	25	U
534-52-1	4,6-Dinitro-2-Methylphenol	25	U
86-30-6	N-Nitrosodiphenylamine (1)	10	U
101-55-3	4-Bromophenyl-phenylether	10	U
118-74-1	Hexachlorobenzene	10	U
87-86-5	Pentachlorophenol	1	J
85-01-8	Phenanthrene	10	U
120-12-7	Anthracene	10	U
86-74-8	Carbazole	10	U
84-74-2	Di-n-Butylphthalate	10	U
206-44-0	Fluoranthene	10	U
129-00-0	Pyrene	1	J
85-68-7	Butylbenzylphthalate	5	BJ
91-94-1	3,3'-Dichlorobenzidine	10	U
56-55-3	Benzo(a)Anthracene	10	U
218-01-9	Chrysene	10	U
117-81-7	bis(2-Ethylhexyl)Phthalate	18	B
117-84-0	Di-n-Octyl Phthalate	10	U
205-99-2	Benzo(b)Fluoranthene	10	U
207-08-9	Benzo(k)Fluoranthene	10	U
50-32-8	Benzo(a)Pyrene	10	U
193-39-5	Indeno(1,2,3-cd)Pyrene	10	U
53-70-3	Dibenz(a,h)Anthracene	10	U
191-24-2	Benzo(g,h,i)Perylene	10	U

(1) - Cannot be separated from Diphenylamine

1B
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

ORIGINAL
CKY32 (MS)

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU

Case No.: 13347

SAS No.: _____

SDG No.: CKY28

Matrix: (soil/water) SOIL

Lab Sample ID: 499456

Sample wt/vol: 30.0 (g/mL) G

Lab File ID: GH099456A15

Level: (low/med) LOW

Date Received: 06/24/92

% Moisture: 18 decanted: (Y/N) N

Date Extracted: 06/26/92

Concentrated Extract Volume: 500.0 (uL)

Date Analyzed: 07/01/92

Injection Volume: 2.0(uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 6.5

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG

Q

CAS NO.	COMPOUND			
108-95-2-----	Phenol	2000		
111-44-4-----	bis(2-Chloroethyl)Ether	400	U	
95-57-8-----	2-Chlorophenol	1700		
541-73-1-----	1,3-Dichlorobenzene	400	U	
106-46-7-----	1,4-Dichlorobenzene	1600		
95-50-1-----	1,2-Dichlorobenzene	400	U	
95-48-7-----	2-Methylphenol	400	U	
108-60-1-----	2,2'-Oxybis(1-Chloropropane)	400	U	
106-44-5-----	4-Methylphenol	400	U	
621-64-7-----	N-Nitroso-Di-n-Propylamine	1400		
67-72-1-----	Hexachloroethane	400	U	
98-95-3-----	Nitrobenzene	400	U	
78-59-1-----	Isophorone	400	U	
88-75-5-----	2-Nitrophenol	400	U	
105-67-9-----	2,4-Dimethylphenol	400	U	
111-91-1-----	bis(2-Chloroethoxy)Methane	400	U	
120-83-2-----	2,4-Dichlorophenol	400	U	
120-82-1-----	1,2,4-Trichlorobenzene	1500		
91-20-3-----	Naphthalene	400	U	
106-47-8-----	4-Chloraniline	400	U	
87-68-3-----	Hexachlorobutadiene	400	U	
59-50-7-----	4-Chloro-3-Methylphenol	1600		
91-57-6-----	2-Methylnaphthalene	400	U	
77-47-4-----	Hexachlorocyclopentadiene	400	U	
88-06-2-----	2,4,6-Trichlorophenol	400	U	
95-95-4-----	2,4,5-Trichlorophenol	980	U	
91-58-7-----	2-Chloronaphthalene	400	U	
88-74-4-----	2-Nitroaniline	980	U	
131-11-3-----	Dimethyl Phthalate	400	U	
208-96-8-----	Acenaphthylene	400	U	
606-20-2-----	2,6-Dinitrotoluene	400	U	
99-09-2-----	3-Nitroaniline	980	U	
83-32-9-----	Acenaphthene	1500		

1C
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

ORIGINAL
CKY32MS

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 18347

SAS No.: _____ SDG No.: CKY28

Matrix: (soil/water) SOIL

Lab Sample ID: 499456

Sample wt/vol: 30.0 (g/mL) G

Lab File ID: GHO99456A15

Level: (low/med) LOW

Date Received: 06/24/92

% Moisture: 18 decanted: (Y/N) N

Date Extracted: 06/26/92

Concentrated Extract Volume: 500.0 (uL)

Date Analyzed: 07/01/92

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 6.5

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG Q

CAS NO.	COMPOUND			
51-28-5-----	2,4-Dinitrophenol	980	U	
100-02-7-----	4-Nitrophenol	3400		
132-64-9-----	Dibenzofuran	400	U	
121-14-2-----	2,4-Dinitrotoluene	1700		
84-66-2-----	Diethylphthalate	400	U	
7005-72-3-----	4-Chlorophenyl-phenylether	400	U	
86-73-7-----	Fluorene	400	U	
100-01-6-----	4-Nitroaniline	980	U	
534-52-1-----	4,6-Dinitro-2-Methylphenol	980	U	
86-30-6-----	N-Nitrosodiphenylamine (1)	400	U	
101-55-3-----	4-Bromophenyl-phenylether	400	U	
118-74-1-----	Hexachlorobenzene	400	U	
87-86-5-----	Pentachlorophenol	1500		
85-01-8-----	Phenanthrene	400	U	
120-12-7-----	Anthracene	400	U	
86-74-8-----	Carbazole	400	U	
84-74-2-----	Di-n-Butylphthalate	400	U	
206-44-0-----	Fluoranthene	400	U	
129-00-0-----	Pyrene	1400		
85-68-7-----	Butylbenzylphthalate	160	BJ	
91-94-1-----	3,3'-Dichlorobenzidine	400	U	
56-55-3-----	Benzo(a)Anthracene	400	U	
218-01-9-----	Chrysene	400	U	
117-81-7-----	bis(2-Ethylhexyl)Phthalate	400	U	
117-84-0-----	Di-n-Octyl Phthalate	400	U	
205-99-2-----	Benzo(b)Fluoranthene	400	U	
207-08-9-----	Benzo(k)Fluoranthene	400	U	
50-32-8-----	Benzo(a)Pyrene	400	U	
193-39-5-----	Indeno(1,2,3-cd)Pyrene	400	U	
53-70-3-----	Dibenz(a,h)Anthracene	400	U	
191-24-2-----	Benzo(g,h,i)Perylene	400	U	

(1) - Cannot be separated from Diphenylamine

1B
SEMICVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CKY32MSD4L
OP/IN
Med

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 18347

SAS No.: _____ SDG No.: CKY28

Matrix: (soil/water) SOIL

Lab Sample ID: 499457

Sample wt/vol: 30.0 (g/mL) G

Lab File ID: GH099457A15

Level: (low/med) LOW

Date Received: 06/24/92

% Moisture: 18 decanted: (Y/N) N

Date Extracted: 06/26/92

Concentrated Extract Volume: 500.0 (uL)

Date Analyzed: 07/01/92

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 6.5

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG

Q

CAS NO.	COMPOUND			
108-95-2	Phenol	2000		
111-44-4	bis(2-Chloroethyl)Ether	400	U	
95-57-8	2-Chlorophenol	1700		
541-73-1	1,3-Dichlorobenzene	400	U	
106-46-7	1,4-Dichlorobenzene	1500		
95-50-1	1,2-Dichlorobenzene	400	U	
95-48-7	2-Methylphenol	400	U	
108-60-1	2,2'-Oxybis(1-Chloropropane)	400	U	
106-44-5	4-Methylphenol	400	U	
621-64-7	N-Nitroso-Di-n-Propylamine	1500		
67-72-1	Hexachloroethane	400	U	
98-95-3	Nitrobenzene	400	U	
78-59-1	Isophorone	400	U	
88-75-5	2-Nitrophenol	400	U	
105-67-9	2,4-Dimethylphenol	400	U	
111-91-1	bis(2-Chloroethoxy)Methane	400	U	
120-83-2	2,4-Dichlorophenol	400	U	
120-82-1	1,2,4-Trichlorobenzene	1500		
91-20-3	Naphthalene	400	U	
106-47-8	4-Chloroaniline	400	U	
87-68-3	Hexachlorobutadiene	400	U	
59-50-7	4-Chloro-3-Methylphenol	1700		
91-57-6	2-Methylnaphthalene	400	U	
77-47-4	Hexachlorocyclopentadiene	400	U	
88-06-2	2,4,6-Trichlorophenol	400	U	
95-95-4	2,4,5-Trichlorophenol	980	U	
91-58-7	2-Chloronaphthalene	400	U	
88-74-4	2-Nitroaniline	980	U	
131-11-3	Dimethyl Phthalate	400	U	
208-96-8	Acenaphthylene	400	U	
606-20-2	2,6-Dinitrotoluene	400	U	
99-09-2	3-Nitroaniline	980	U	
83-32-9	Acenaphthene	1600		

1C
SEMICVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

ORIGINAL
(Red)

CKY32MSD

Lab Name: COMPUCHEM.RTPContract: 68D10083Lab Code: COMPUCase No.: 18347

SAS No.: _____

SDG No.: CKY28Matrix: (soil/water) SOILLab Sample ID: 499457Sample wt/vol: 30.0 (g/mL) GLab File ID: GH099457A15Level: (low/med) LOWDate Received: 06/24/92% Moisture: 18 decanted: (Y/N) NDate Extracted: 06/26/92Concentrated Extract Volume: 500.0 (uL)Date Analyzed: 07/01/92Injection Volume: 2.0(uL)Dilution Factor: 1.0GPC Cleanup: (Y/N) Y pH: 6.5CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG Q

CAS NO.	COMPCUND	Q
51-28-5-----	2,4-Dinitrophenol	980 U
100-02-7-----	4-Nitrophenol	3600 U
132-64-9-----	Dibenzofuran	400 U
121-14-2-----	2,4-Dinitrotoluene	1800 U
84-66-2-----	Diethylphthalate	400 U
7005-72-3-----	4-Chlorophenyl-phenylether	400 U
86-73-7-----	Fluorene	400 U
100-01-6-----	4-Nitroaniline	980 U
534-52-1-----	4,6-Dinitro-2-Methylphenol	980 U
86-30-6-----	N-Nitrosodiphenylamine (1)	400 U
101-55-3-----	4-Bromophenyl-phenylether	400 U
118-74-1-----	Hexachlorobenzene	400 U
87-86-5-----	Pentachlorophenol	1500 U
85-01-8-----	Phenanthrene	400 U
120-12-7-----	Anthracene	400 U
86-74-8-----	Carbazole	400 U
84-74-2-----	Di-n-Butylphthalate	400 U
206-44-0-----	Fluoranthene	400 U
129-00-0-----	Pyrene	1500 U
65-68-7-----	Butylbenzylphthalate	150 BJ
91-94-1-----	3,3'-Dichlorobenzidine	400 U
56-55-3-----	Benzo(a)Anthracene	400 U
218-01-9-----	Chrysene	400 U
117-81-7-----	bis(2-Ethylhexyl)Phthalate	400 U
117-84-0-----	Di-n-Octyl Phthalate	400 U
205-99-2-----	Benzo(b)Fluoranthene	48 JX
207-08-9-----	Benzo(k)Fluoranthene	48 JX
50-32-8-----	Benzo(a)Pyrene	400 U
193-39-5-----	Indeno(1,2,3-cd)Pyrene	400 U
53-70-3-----	Dibenz(a,h)Anthracene	400 U
191-24-2-----	Benzo(g,h,i)Perylene	400 U

(1) - Cannot be separated from Diphenylamine

1D
PESTICIDE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

Original

Mod

CKY18MS

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 18347 SAS No.:

SDG No.: CKY14

Matrix: (soil/water) WATER

Lab Sample ID: 499432

Sample wt/vol: 500(g/ml)ML

Lab File ID:

% Moisture: decanted: (Y/N)

Date Received: 06/24/92

Extraction: (SepF/Cont/Sonc) SEPF

Date Extracted: 06/26/92

Concentrated Extract Volume: 5000(uL)

Date Analyzed: 06/30/92

Injection Volume: 2.0(uL)

Dilution Factor: 1

GPC Cleanup: (Y/N) N pH:

Sulfur Cleanup: (Y/N) N

CONCENTRATION UNITS:

(ug/L or ug/Kg) UG/L

Q

319-84-6-----	alpha-BHC	0.050	U
319-85-7-----	beta-BHC	0.050	U
319-86-8-----	delta-BHC	0.050	U
58-89-9-----	gamma-BHC (Lindane)	0.44	
76-44-8-----	Heptachlor	0.40	
309-00-2-----	Aldrin	0.39	
1024-57-3-----	Heptachlor epoxide	0.050	U
959-98-8-----	Endosulfan I	0.050	U
60-57-1-----	Dieldrin	0.76	
72-55-9-----	4,4'-DDE	0.10	U
72-20-8-----	Endrin	0.88	
33213-65-9-----	Endosulfan II	0.10	U
72-54-8-----	4,4'-DDD	0.10	U
1031-07-8-----	Endosulfan sulfate	0.10	U
50-29-3-----	4,4'-DDT	0.92	
72-43-5-----	Methoxychlor	0.017	JP
53494-70-5-----	Endrin ketone	0.0099	J
7421-93-4-----	Endrin aldehyde	0.10	U
5103-71-9-----	alpha-Chlordane	0.050	U
5103-74-2-----	gamma-Chlordane	0.050	U
8001-35-2-----	Toxaphene	5.0	U
12674-11-2-----	Aroclor-1016	1.0	U
11104-28-2-----	Aroclor-1221	2.0	U
11141-16-5-----	Aroclor-1232	1.0	U
53469-21-9-----	Aroclor-1242	1.0	U
12672-29-6-----	Aroclor-1248	1.0	U
11097-69-1-----	Aroclor-1254	1.0	U
11096-82-5-----	Aroclor-1260	1.0	U

ID
PESTICIDE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

ORIGINAL

(Red)
CKY18MSD

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 18347 SAS No.:

SDG No.: CKY14

Matrix: (soil/water) WATER

Lab Sample ID: 499433

Sample wt/vol: 500(g/ml) ML

Lab File ID:

Moisture: decanted: (Y/N)

Date Received: 06/24/92

Extraction: (SepF/Cont/Sonc) SEPF

Date Extracted: 06/26/92

Concentrated Extract Volume: 5000(uL)

Date Analyzed: 06/30/92

Injection Volume: 2.0(uL)

Dilution Factor: 1

GPC Cleanup: (Y/N) N

pH:

Sulfur Cleanup: (Y/N) N

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/L

Q

319-84-6-----alpha-BHC	0.050	U
319-85-7-----beta-BHC	0.050	U
319-86-8-----delta-BHC	0.050	U
58-89-9-----gamma-BHC (Lindane)	0.41	
76-44-8-----Heptachlor	0.38	
309-00-2-----Aldrin	0.37	
1024-57-3-----Heptachlor epoxide	0.050	U
959-98-8-----Endosulfan I	0.050	U
60-57-1-----Dieldrin	0.72	
72-55-9-----4,4'-DDE	0.10	U
72-20-8-----Endrin	0.84	
33213-65-9-----Endosulfan II	0.10	U
72-54-8-----4,4'-DDD	0.10	U
1031-07-8-----Endosulfan sulfate	0.10	U
50-29-3-----4,4'-DDT	0.88	
72-43-5-----Methoxychlor	0.013	JP
53494-70-5-----Endrin ketone	0.0074	J
7421-93-4-----Endrin aldehyde	0.10	U
5103-71-9-----alpha-Chlordane	0.050	U
5103-74-2-----gamma-Chlordane	0.050	U
8001-35-2-----Toxaphene	5.0	U
12674-11-2-----Aroclor-1016	1.0	U
11104-28-2-----Aroclor-1221	2.0	U
11141-16-5-----Aroclor-1232	1.0	U
53469-21-9-----Aroclor-1242	1.0	U
12672-29-6-----Aroclor-1248	1.0	U
11097-69-1-----Aroclor-1254	1.0	U
11096-82-5-----Aroclor-1260	1.0	U

1D
PESTICIDE ORGANICS ANALYSIS DATA SHEET

EPA/SAMPLE NO.
(Ref)

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

CKY32MS

Lab Code: COMPU Case No.: 18347 SAS No.:

SDG No.: CKY28

Matrix: (soil/water) SOIL

Lab Sample ID: 499460

Sample wt/vol: 30.00(g/ml)G

Lab File ID:

% Moisture: 18 decanted: (Y/N) N

Date Received: 06/24/92

Extraction: (Sep/Cont/Sonic) SONC

Date Extracted: 06/26/92

Concentrated Extract Volume: 5000(uL)

Date Analyzed: 07/06/92

Injection Volume: 2.0(uL)

Dilution Factor: 1

GPC Cleanup: (Y/N) Y pH: 6.5

Sulfur Cleanup: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) UG/KG	Q
319-84-6-----	alpha-BHC	2.1	U
319-85-7-----	beta-BHC	2.1	U
319-86-8-----	delta-BHC	0.093	JP
58-89-9-----	gamma-BHC (Lindane)	13	
76-44-8-----	Heptachlor	11	
309-00-2-----	Aldrin	12	
1024-57-3-----	Heptachlor epoxide	0.21	J
959-98-8-----	Endosulfan I	2.1	U
60-57-1-----	Dieldrin	25	
72-55-9-----	4,4'-DDE	4.0	U
72-20-8-----	Endrin	27	
33213-65-9-----	Endosulfan II	4.0	U
72-54-8-----	4,4'-DDD	4.0	U
1031-07-8-----	Endosulfan sulfate	4.0	U
50-29-3-----	4,4'-DDT	25	P
72-43-5-----	Methoxychlor	21	U
53494-70-5-----	Endrin ketone	4.0	U
7421-93-4-----	Endrin aldehyde	4.0	U
5103-71-9-----	alpha-Chlordane	2.1	U
5103-74-2-----	gamma-Chlordane	2.1	U
8001-35-2-----	Toxaphene	210	U
12674-11-2-----	Aroclor-1016	40	U
11104-28-2-----	Aroclor-1221	82	U
11141-16-5-----	Aroclor-1232	40	U
53469-21-9-----	Aroclor-1242	40	U
12672-29-6-----	Aroclor-1248	40	U
11097-69-1-----	Aroclor-1254	40	U
11096-82-5-----	Aroclor-1260	40	U

1D
PESTICIDE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

CKY32MSD

Lab Code: COMPU Case No.: 18347 SAS No.:

SDG No.: CKY28

Matrix: (soil/water) SOIL

Lab Sample ID: 499461

Sample wt/vol: 30.00(g/ml)G

Lab File ID:

% Moisture: 13 decanted: (Y/N) N

Date Received: 06/24/92

Extraction: (SepF/Cont/Sonc) SONC

Date Extracted: 06/26/92

Concentrated Extract Volume: 5000(uL)

Date Analyzed: 07/06/92

Injection Volume: 2.0(uL)

Dilution Factor: 1

GPC Cleanup: (Y/N) Y pH: 6.5

Sulfur Cleanup: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/KG</u>	Q
319-84-5-----	alpha-BHC	2.1	U
319-85-7-----	beta-BHC	2.1	U
319-86-8-----	delta-BHC	0.22	JP
58-89-9-----	gamma-BHC (Lindane)	13	
76-44-8-----	Heptachlor	12	
309-00-2-----	Aldrin	13	
1024-57-3-----	Heptachlor epoxide	2.1	U
959-98-8-----	Endosulfan I	2.1	U
60-57-1-----	Dieldrin	27	
72-55-9-----	1,4'-DDE	4.0	U
72-20-8-----	Endrin	28	
33213-65-9-----	Endosulfan II	4.0	U
72-54-8-----	1,4'-DDD	4.0	U
1031-07-8-----	Endosulfan sulfate	4.0	U
50-29-3-----	1,4'-DDT	26	
72-43-5-----	Methoxychlor	21	U
53494-70-5-----	Endrin ketone	1.2	J
7421-93-4-----	Endrin aldehyde	4.0	U
5103-71-9-----	alpha-Chlordane	2.1	U
5103-74-2-----	gamma-Chlordane	2.1	U
8001-35-2-----	Toxaphene	210	U
12674-11-2-----	Aroclor-1016	40	U
11104-28-2-----	Aroclor-1221	82	U
11141-16-5-----	Aroclor-1232	40	U
53469-21-9-----	Aroclor-1242	40	U
12672-29-6-----	Aroclor-1248	40	U
11097-69-1-----	Aroclor-1254	40	U
11096-82-5-----	Aroclor-1260	40	U

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

O/R/05

(Red)

VBLKC7

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 18347

SAS No.: _____ SDG No.: CKY28

Matrix: (soil/water) SOIL

Lab Sample ID: VBLKC7

Sample wt/vol: 5.0 (g/mL) G

Lab File ID: GH099517B54

Level: (low/med) LOW

Date Received: _____

% Moisture: not dec. _____

Date Analyzed: 06/24/92

Column: DB-624 ID: 0.530 (mm)

Dilution Factor: 1.0

Soil Extract Volume: _____ (uL)

Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

(ug/L or ug/Kg) UG/KG

Q

CAS NO.	COMPOUND	10	U
74-87-3-----	Chloromethane	10	U
74-83-9-----	Bromomethane	10	U
75-01-4-----	Vinyl Chloride	10	U
75-00-3-----	Chloroethane	10	U
75-09-2-----	Methylene Chloride	10	
67-64-1-----	Acetone	28	
75-15-0-----	Carbon Disulfide	10	U
75-35-4-----	1,1-Dichloroethene	10	U
75-34-3-----	1,1-Dichloroethane	10	U
540-59-0-----	1,2-Dichloroethene (total)	10	U
67-66-3-----	Chloroform	10	U
107-06-2-----	1,2-Dichloroethane	10	U
78-93-3-----	2-Butanone	10	U
71-55-6-----	1,1,1-Trichloroethane	10	U
56-23-5-----	Carbon Tetrachloride	10	U
75-27-4-----	Bromodichloromethane	10	U
78-87-5-----	1,2-Dichloropropane	10	U
10061-01-5-----	cis-1,3-Dichloropropene	10	U
79-01-6-----	Trichloroethene	10	U
124-48-1-----	Dibromochloromethane	10	U
79-00-5-----	1,1,2-Trichloroethane	10	U
71-43-2-----	Benzene	10	U
10061-02-6-----	Trans-1,3-Dichloropropene	10	U
75-25-2-----	Bromoform	10	U
108-10-1-----	4-Methyl-2-Pentanone	10	U
591-78-6-----	2-Hexanone	10	U
127-18-4-----	Tetrachloroethene	10	U
79-34-5-----	1,1,2,2-Tetrachloroethane	10	U
108-88-3-----	Toluene	10	U
108-90-7-----	Chlorobenzene	10	U
100-41-4-----	Ethylbenzene	10	U
100-42-5-----	Styrene	10	U
1330-20-7-----	Xylene (total)	10	U

1E
 VOLATILE ORGANICS ANALYSIS DATA SHEET
 TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

(Read)

VBLKC7

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 18347

SAS No.: _____ SDG No.: CKY28

Matrix: (soil/water) SOIL

Lab Sample ID: VBLKC7

Sample wt/vol: 5.0 (g/mL) G

Lab File ID: GH099517B54

Level: (low/med) LOW

Date Received: _____

% Moisture: not dec. _____

Date Analyzed: 06/24/92

GC Column: DB-624 ID: 0.530 (mm)

Dilution Factor: 1.0

Soil Extract Volume: _____ (uL)

Soil Aliquot Volume: _____ (uL)

Number TICs found: 1

CONCENTRATION UNITS:

(ug/L or ug/Kg) UG/KG

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1.	LABORATORY ARTIFACT	2.18	36	J-Hfa 25/9

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

(R00)

VBLKC8

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 18347

SAS No.: _____ SDG No.: CKY28

Matrix: (soil/water) SOIL

Lab Sample ID: VBLKC8

Sample wt/vol: 5.0 (g/mL) G

Lab File ID: GH099518C54

Level: (low/med) LOW

Date Received: _____

% Moisture: not dec. _____

Date Analyzed: 06/25/92

GC Column: DB-624 ID: 0.530 (mm)

Dilution Factor: 1.0

Soil Extract Volume: _____ (uL)

Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

(ug/L or ug/Kg) UG/KG

Q

CAS NO.	COMPOUND		
74-87-3-----	Chloromethane	10	U
74-83-9-----	Bromomethane	10	U
75-01-4-----	Vinyl Chloride	10	U
75-00-3-----	Chloroethane	10	U
75-09-2-----	Methylene Chloride	30	
67-64-1-----	Acetone	31	
75-15-0-----	Carbon Disulfide	10	U
75-35-4-----	1,1-Dichloroethene	10	U
75-34-3-----	1,1-Dichloroethane	10	U
540-59-0-----	1,2-Dichloroethene (total)	10	U
67-66-3-----	Chloroform	10	U
107-06-2-----	1,2-Dichloroethane	10	U
78-93-3-----	2-Butanone	10	U
71-55-6-----	1,1,1-Trichloroethane	10	U
56-23-5-----	Carbon Tetrachloride	10	U
75-27-4-----	Bromodichloromethane	10	U
78-87-5-----	1,2-Dichloropropane	10	U
10061-01-5-----	cis-1,3-Dichloropropene	10	U
79-01-6-----	Trichloroethene	10	U
124-48-1-----	Dibromochloromethane	10	U
79-00-5-----	1,1,2-Trichloroethane	10	U
71-43-2-----	Benzene	10	U
10061-02-6-----	Trans-1,3-Dichloropropene	10	U
75-25-2-----	Bromoform	10	U
108-10-1-----	4-Methyl-2-Pentanone	10	U
591-78-6-----	2-Hexanone	10	U
127-18-4-----	Tetrachloroethene	10	U
79-34-5-----	1,1,2,2-Tetrachloroethane	10	U
108-88-3-----	Toluene	10	U
108-90-7-----	Chlorobenzene	10	U
100-41-4-----	Ethylbenzene	10	U
100-42-5-----	Styrene	10	U
1330-20-7-----	Xylene (total)	10	U

1E
VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

ORIGINAL

(Rev)

VBLKC8

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 18347

SAS No.: _____ SDG No.: CKY28

Matrix: (soil/water) SOIL

Lab Sample ID: VBLKC8

Sample wt/vol: 5.0 (g/mL) G

Lab File ID: GH099518C54

Level: (low/med) LOW

Date Received: _____

% Moisture: not dec. _____

Date Analyzed: 06/25/92

Column: DB-624 ID: 0.530 (mm)

Dilution Factor: 1.0

Soil Extract Volume: _____ (uL)

Soil Aliquot Volume: _____ (uL)

Number TICs found: 1

CONCENTRATION UNITS:

(ug/L or ug/Kg) UG/KG

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1.	LABORATORY ARTIFACT	2.20	28	BJ 18/9 28

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

09/15/92

(Ref'd)

VBLKKX

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 18347

SAS No.: _____ SDG No.: CKY14

Matrix: (soil/water) WATER

Lab Sample ID: VBLKKX

Sample wt/vol: 5.0 (g/mL) ML

Lab File ID: CB920624A56

Level: (low/med) LOW

Date Received: _____

% Moisture: not dec. _____

Date Analyzed: 06/24/92

Column: DB624 ID: 0.530 (mm)

Dilution Factor: 1.0

Soil Extract Volume: _____ (uL)

Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

(ug/L or ug/Kg) UG/L

Q

CAS NO.	COMPOUND	Q
74-87-3-----	Chloromethane	10 U
74-83-9-----	Bromomethane	10 U
75-01-4-----	Vinyl Chloride	10 U
75-00-3-----	Chloroethane	10 U
75-09-2-----	Methylene Chloride	3 J
67-64-1-----	Acetone	13
75-15-0-----	Carbon Disulfide	10 U
75-35-4-----	1,1-Dichloroethene	10 U
75-34-3-----	1,1-Dichloroethane	10 U
540-59-0-----	1,2-Dichloroethene (total)	10 U
67-66-3-----	Chloroform	10 U
107-06-2-----	1,2-Dichloroethane	10 U
78-93-3-----	2-Butanone	10 U
71-55-6-----	1,1,1-Trichloroethane	10 U
56-23-5-----	Carbon Tetrachloride	10 U
75-27-4-----	Bromodichloromethane	10 U
78-87-5-----	1,2-Dichloropropane	10 U
10061-01-5-----	cis-1,3-Dichloropropene	10 U
79-01-6-----	Trichloroethene	10 U
124-48-1-----	Dibromochloromethane	10 U
79-00-5-----	1,1,2-Trichloroethane	10 U
71-43-2-----	Benzene	10 U
10061-02-6-----	Trans-1,3-Dichloropropene	10 U
75-25-2-----	Bromoform	10 U
108-10-1-----	4-Methyl-2-Pentanone	10 U
591-78-6-----	2-Hexanone	10 U
127-18-4-----	Tetrachloroethene	10 U
79-34-5-----	1,1,2,2-Tetrachloroethane	10 U
108-88-3-----	Toluene	10 U
108-90-7-----	Chlorobenzene	10 U
100-41-4-----	Ethylbenzene	10 U
100-42-5-----	Styrene	10 U
1330-20-7-----	Xylene (total)	10 U

1E
VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

ORIGINAL
Med

VBLKKX

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 18347

SAS No.: _____ SDG No.: CKY14

Matrix: (soil/water) WATER

Lab Sample ID: VBLKKX

Sample wt/vol: 5.0 (g/mL) ML

Lab File ID: CB920624A56

Level: (low/med) LOW

Date Received: _____

% Moisture: not dec. _____

Date Analyzed: 06/24/92

GC Column: DB624 ID: 0.530 (mm)

Dilution Factor: 1.0

Soil Extract Volume: _____ (uL)

Soil Aliquot Volume: _____ (uL)

Number TICs found: 0

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
=====	=====	=====	=====	=====

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

(low/med)
VBLKKA

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 18347

SAS No.: _____ SDG No.: CKY14

Matrix: (soil/water) WATER

Lab Sample ID: VBLKKA

Sample wt/vol: 5.0 (g/mL) ML

Lab File ID: CD920625A56

Level: (low/med) LOW

Date Received: _____

% Moisture: not dec. _____

Date Analyzed: 06/25/92

GC Column: DB624 ID: 0.530 (mm)

Dilution Factor: 1.0

Soil Extract Volume: _____ (uL)

Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

(ug/L or ug/Kg) UG/L

Q

<u>74-87-3-----Chloromethane</u>	<u>10</u>	<u>U</u>
<u>74-83-9-----Bromomethane</u>	<u>10</u>	<u>U</u>
<u>75-01-4-----Vinyl Chloride</u>	<u>10</u>	<u>U</u>
<u>75-00-3-----Chloroethane</u>	<u>10</u>	<u>U</u>
<u>75-09-2-----Methylene Chloride</u>	<u>3</u>	<u>J</u>
<u>67-64-1-----Acetone</u>	<u>13</u>	
<u>75-15-0-----Carbon Disulfide</u>	<u>10</u>	<u>U</u>
<u>75-35-4-----1,1-Dichloroethene</u>	<u>10</u>	<u>U</u>
<u>75-34-3-----1,1-Dichloroethane</u>	<u>10</u>	<u>U</u>
<u>540-59-0-----1,2-Dichloroethene (total)</u>	<u>10</u>	<u>U</u>
<u>67-66-3-----Chloroform</u>	<u>10</u>	<u>U</u>
<u>107-06-2-----1,2-Dichloroethane</u>	<u>10</u>	<u>U</u>
<u>78-93-3-----2-Butanone</u>	<u>10</u>	<u>U</u>
<u>71-55-6-----1,1,1-Trichloroethane</u>	<u>10</u>	<u>U</u>
<u>56-23-5-----Carbon Tetrachloride</u>	<u>10</u>	<u>U</u>
<u>75-27-4-----Bromodichloromethane</u>	<u>10</u>	<u>U</u>
<u>78-87-5-----1,2-Dichloropropane</u>	<u>10</u>	<u>U</u>
<u>10061-01-5-----cis-1,3-Dichloropropene</u>	<u>10</u>	<u>U</u>
<u>79-01-6-----Trichloroethene</u>	<u>10</u>	<u>U</u>
<u>124-48-1-----Dibromochloromethane</u>	<u>10</u>	<u>U</u>
<u>79-00-5-----1,1,2-Trichloroethane</u>	<u>10</u>	<u>U</u>
<u>71-43-2-----Benzene</u>	<u>10</u>	<u>U</u>
<u>10061-02-6-----Trans-1,3-Dichloropropene</u>	<u>10</u>	<u>U</u>
<u>75-25-2-----Bromoform</u>	<u>10</u>	<u>U</u>
<u>108-10-1-----4-Methyl-2-Pentanone</u>	<u>10</u>	<u>U</u>
<u>591-78-6-----2-Hexanone</u>	<u>10</u>	<u>U</u>
<u>127-18-4-----Tetrachloroethene</u>	<u>10</u>	<u>U</u>
<u>79-34-5-----1,1,2,2-Tetrachloroethane</u>	<u>10</u>	<u>U</u>
<u>108-88-3-----Toluene</u>	<u>10</u>	<u>U</u>
<u>108-90-7-----Chlorobenzene</u>	<u>10</u>	<u>U</u>
<u>100-41-4-----Ethylbenzene</u>	<u>10</u>	<u>U</u>
<u>100-42-5-----Styrene</u>	<u>10</u>	<u>U</u>
<u>1330-20-7-----Xylene (total)</u>	<u>10</u>	<u>U</u>

1E
VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

ORIGINAL

(Regd)

VBLKKA

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 18347

SAS No.: _____ SDG No.: CKY14

Matrix: (soil/water) WATER

Lab Sample ID: VBLKKA

Sample wt/vol: 5.0 (g/mL) ML

Lab File ID: CD920625A56

Level: (low/med) LOW

Date Received: _____

% Moisture: not dec. _____

Date Analyzed: 06/25/92

GC Column: DB624 ID: 0.530 (mm)

Dilution Factor: 1.0

Soil Extract Volume: _____ (uL)

Soil Aliquot Volume: _____ (uL)

Number TICs found: 0

CONCENTRATION UNITS:

(ug/L or ug/Kg) UG/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
=====	=====	=====	=====	=====

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

VBLKVT^{ORIGINAL}
(Red)

Lab Code: COMPU Case No.: 18347

SAS No.: _____ SDG No.: CKY14

Matrix: (soil/water) WATER

Lab Sample ID: VBLKVT

Sample wt/vol: 5.0 (g/mL) ML

Lab File ID: CB920625B56

Level: (low/med) LOW

Date Received: _____

% Moisture: not dec. _____

Date Analyzed: 06/25/92

GC Column: DB624 ID: 0.530 (mm)

Dilution Factor: 1.0

Soil Extract Volume: _____ (uL)

Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

(ug/L or ug/Kg) UG/L

Q

74-87-3-----Chloromethane	10	U
74-83-9-----Bromomethane	10	U
75-01-4-----Vinyl Chloride	10	U
75-00-3-----Chloroethane	10	U
75-09-2-----Methylene Chloride	8	J
67-64-1-----Acetone	5	J
75-15-0-----Carbon Disulfide	10	U
75-35-4-----1,1-Dichloroethene	10	U
75-34-3-----1,1-Dichloroethane	10	U
540-59-0-----1,2-Dichloroethene (total)	10	U
67-66-3-----Chloroform	10	U
107-06-2-----1,2-Dichloroethane	10	U
78-93-3-----2-Butanone	10	U
71-55-6-----1,1,1-Trichloroethane	10	U
56-23-5-----Carbon Tetrachloride	10	U
75-27-4-----Bromodichloromethane	10	U
78-87-5-----1,2-Dichloropropane	10	U
10061-01-5-----cis-1,3-Dichloropropene	10	U
79-01-6-----Trichloroethene	10	U
124-48-1-----Dibromoethane	10	U
79-00-5-----1,1,2-Trichloroethane	10	U
71-43-2-----Benzene	10	U
10061-02-6-----Trans-1,3-Dichloropropene	10	U
75-25-2-----Bromoform	10	U
108-10-1-----4-Methyl-2-Pentanone	10	U
591-78-6-----2-Hexanone	10	U
127-18-4-----Tetrachloroethene	10	U
79-34-5-----1,1,2,2-Tetrachloroethane	10	U
108-88-3-----Toluene	10	U
108-90-7-----Chlorobenzene	10	U
100-41-4-----Ethylbenzene	10	U
100-42-5-----Styrene	10	U
1330-20-7-----Xylene (total)	10	U

1E
VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.
(Recd)

VBLKVT

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 18347

SAS No.: _____ SDG No.: CKY14

Matrix: (soil/water) WATER

Lab Sample ID: VBLKVT

Sample wt/vol: 5.0 (g/mL) ML

Lab File ID: CB920625B56

Level: (low/med) LOW

Date Received: _____

% Moisture: not dec.

Date Analyzed: 06/25/92

Column: DB624 ID: 0.530 (mm)

Dilution Factor: 1.0

Soil Extract Volume: _____ (uL)

Soil Aliquot Volume: _____ (uL)

Number TICs found: 0

CONCENTRATION UNITS:

(ug/L or ug/Kg) UG/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
=====	=====	=====	=====	=====

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA/SAMPLE NO.
(Red)

VBLKD3

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU

Case No.: 13347

SAS No.: _____

SDG No.: CKY28

Matrix: (soil/water) SOIL

Lab Sample ID: VBLKD3

Sample wt/vol: 5.0 (g/mL) G

Lab File ID: GH099662A54

Level: (low/med) LOW

Date Received: _____

% Moisture: not dec. _____

Date Analyzed: 06/25/92

GC Column: DB-624 ID: 0.530 (mm)

Dilution Factor: 1.0

Soil Extract Volume: _____ (uL)

Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

(ug/L or ug/Kg) UG/KG

Q

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/KG</u>	Q
74-87-3-----	Chloromethane	10	U
74-83-9-----	Bromomethane	10	U
75-01-4-----	Vinyl Chloride	10	U
75-00-3-----	Chloroethane	10	U
75-09-2-----	Methylene Chloride	30	
67-64-1-----	Acetone	21	
75-15-0-----	Carbon Disulfide	10	U
75-35-4-----	1,1-Dichloroethene	10	U
75-34-3-----	1,1-Dichloroethane	10	U
540-59-0-----	1,2-Dichloroethene (total)	10	U
67-66-3-----	Chloroform	10	U
107-06-2-----	1,2-Dichloroethane	10	U
78-93-3-----	2-Butanone	10	U
71-55-6-----	1,1,1-Trichloroethane	10	U
56-23-5-----	Carbon Tetrachloride	10	U
75-27-4-----	Bromodichloromethane	10	U
78-87-5-----	1,2-Dichloropropane	10	U
10061-01-5-----	cis-1,3-Dichloropropene	10	U
79-01-6-----	Trichloroethene	10	U
124-48-1-----	Dibromochloromethane	10	U
79-00-5-----	1,1,2-Trichloroethane	10	U
71-43-2-----	Benzene	10	U
10061-02-6-----	Trans-1,3-Dichloropropene	10	U
75-25-2-----	Bromoform	10	U
108-10-1-----	4-Methyl-2-Pentanone	10	U
591-78-6-----	2-Hexanone	10	U
127-18-4-----	Tetrachloroethene	10	U
79-34-5-----	1,1,2,2-Tetrachloroethane	10	U
108-88-3-----	Toluene	10	U
108-90-7-----	Chlorobenzene	10	U
100-41-4-----	Ethylbenzene	10	U
100-42-5-----	Styrene	10	U
1330-20-7-----	Xylene (total)	10	U

1E
VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

VBLKD3

ORIGINAL
(Red)

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 18347

SAS No.: _____ SDG No.: CKY28

Matrix: (soil/water) SOIL

Lab Sample ID: VBLKD3

Sample wt/vol: 5.0 (g/mL) G

Lab File ID: GH099662A54

Level: (low/med) LOW

Date Received: _____

% Moisture: not dec. _____

Date Analyzed: 06/25/92

GC Column: DB-624 ID: 0.530 (mm)

Dilution Factor: 1.0

Soil Extract Volume: _____ (uL)

Soil Aliquot Volume: _____ (uL)

Number TICs found: 1

CONCENTRATION UNITS:

(ug/L or ug/Kg) UG/KG

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1.	LABORATORY ARTIFACT	2.17	20	3.18912512

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

ORIGINAL
(Red)

EPA SAMPLE NO.

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

VBLKD9

Lab Code: COMPU Case No.: 18347

SAS No.: _____

SDG No.: CKY28

Matrix: (soil/water) SOIL

Lab Sample ID: VBLKD9

Sample wt/vol: 5.0 (g/mL) G

Lab File ID: GH099886A54

Level: (low/med) LOW

Date Received: _____

% Moisture: not dec. _____

Date Analyzed: 06/26/92

GC Column: DB-624 ID: 0.530 (mm)

Dilution Factor: 1.0

Soil Extract Volume: _____ (uL)

Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

(ug/L or ug/Kg) UG/KG

Q

<u>74-87-3-----Chloromethane</u>	<u>10</u>	<u>U</u>
<u>74-83-9-----Bromomethane</u>	<u>10</u>	<u>U</u>
<u>75-01-4-----Vinyl Chloride</u>	<u>10</u>	<u>U</u>
<u>75-00-3-----Chloroethane</u>	<u>10</u>	<u>U</u>
<u>75-09-2-----Methylene Chloride</u>	<u>23</u>	
<u>67-64-1-----Acetone</u>	<u>18</u>	
<u>75-15-0-----Carbon Disulfide</u>	<u>10</u>	<u>U</u>
<u>75-35-4-----1,1-Dichloroethene</u>	<u>10</u>	<u>U</u>
<u>75-34-3-----1,1-Dichloroethane</u>	<u>10</u>	<u>U</u>
<u>540-59-0-----1,2-Dichloroethene (total)</u>	<u>10</u>	<u>U</u>
<u>67-66-3-----Chloroform</u>	<u>10</u>	<u>U</u>
<u>107-06-2-----1,2-Dichloroethane</u>	<u>10</u>	<u>U</u>
<u>78-93-3-----2-Butanone</u>	<u>10</u>	<u>U</u>
<u>71-55-6-----1,1,1-Trichloroethane</u>	<u>10</u>	<u>U</u>
<u>56-23-5-----Carbon Tetrachloride</u>	<u>10</u>	<u>U</u>
<u>75-27-4-----Bromodichloromethane</u>	<u>10</u>	<u>U</u>
<u>78-37-5-----1,2-Dichloropropane</u>	<u>10</u>	<u>U</u>
<u>10061-01-5-----cis-1,3-Dichloropropene</u>	<u>10</u>	<u>U</u>
<u>79-01-6-----Trichloroethene</u>	<u>10</u>	<u>U</u>
<u>124-48-1-----Dibromochloromethane</u>	<u>10</u>	<u>U</u>
<u>79-00-5-----1,1,2-Trichloroethane</u>	<u>10</u>	<u>U</u>
<u>71-43-2-----Benzene</u>	<u>10</u>	<u>U</u>
<u>10061-02-6-----Trans-1,3-Dichloropropene</u>	<u>10</u>	<u>U</u>
<u>75-25-2-----Bromoform</u>	<u>10</u>	<u>U</u>
<u>108-10-1-----4-Methyl-2-Pentanone</u>	<u>10</u>	<u>U</u>
<u>591-78-6-----2-Hexanone</u>	<u>10</u>	<u>U</u>
<u>127-18-4-----Tetrachloroethene</u>	<u>10</u>	<u>U</u>
<u>79-34-5-----1,1,2,2-Tetrachloroethane</u>	<u>10</u>	<u>U</u>
<u>108-88-3-----Toluene</u>	<u>10</u>	<u>U</u>
<u>108-90-7-----Chlorobenzene</u>	<u>10</u>	<u>U</u>
<u>100-41-4-----Ethylbenzene</u>	<u>10</u>	<u>U</u>
<u>100-42-5-----Styrene</u>	<u>10</u>	<u>U</u>
<u>1330-20-7-----Xylene (total)</u>	<u>10</u>	<u>U</u>

1E
VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

ORIGINAL
fmed
VBLKD9

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 18347 SAS No.: _____ SDG No.: CKY28

Matrix: (soil/water) SOIL

Lab Sample ID: VBLKD9

Sample wt/vol: 5.0 (g/mL) G

Lab File ID: GH099886A54

Level: (low/med) LOW

Date Received: _____

% Moisture: not dec. _____

Date Analyzed: 06/26/92

GC Column: DB-624 ID: 0.530 (mm)

Dilution Factor: 1.0

Soil Extract Volume: _____ (uL)

Soil Aliquot Volume: _____ (uL)

Number TICs found: 1

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1.	LABORATORY ARTIFACT	2.15	19	J-H 9/23/92

1B
SEMICVOLATILE ORGANICS ANALYSIS DATA SHEET

ORIG/EPA SAMPLE NO.

(Recd)

SBLK13

Lab Name: COMPUCHEM, RTPContract: 68D10083Lab Code: COMPU Case No.: 18347SAS No.: _____ SDG No.: CKY14Matrix: (soil/water) WATERLab Sample ID: SBLK13Sample wt/vol: 1000 (g/mL) MLLab File ID: GH099898A07Level: (low/med) LOW

Date Received: _____

% Moisture: _____ decanted: (Y/N) _____

Date Extracted: 06/25/92Concentrated Extract Volume: 1000 (uL)Date Analyzed: 06/29/92Injection Volume: 2.0 (uL)Dilution Factor: 1.0GPC Cleanup: (Y/N) N pH: _____

CONCENTRATION UNITS:

CAS NO. COMPOUND (ug/L or ug/Kg) UG/L Q

108-95-2-----	Phenol	1	J
111-44-4-----	bis(2-Chloroethyl)Ether	10	U
95-57-8-----	2-Chlorophenol	10	U
541-73-1-----	1,3-Dichlorobenzene	10	U
106-46-7-----	1,4-Dichlorobenzene	10	U
95-50-1-----	1,2-Dichlorobenzene	10	U
95-48-7-----	2-Methylphenol	10	U
108-60-1-----	2,2'-Oxybis(1-Chloropropane)	10	U
106-44-5-----	4-Methylphenol	10	U
621-64-7-----	N-Nitroso-Di-n-Propylamine	10	U
67-72-1-----	Hexachloroethane	10	U
98-95-3-----	Nitrobenzene	10	U
78-59-1-----	Iscophonone	10	U
88-75-5-----	2-Nitrophenol	10	U
105-67-9-----	2,4-Dimethylphenol	10	U
111-91-1-----	bis(2-Chloroethoxy)Methane	10	U
120-83-2-----	2,4-Dichlorophenol	10	U
120-82-1-----	1,2,4-Trichlorobenzene	10	U
91-20-3-----	Naphthalene	10	U
106-47-8-----	4-Chloroaniline	10	U
87-68-3-----	Hexachlorobutadiene	10	U
59-50-7-----	4-Chloro-3-Methylphenol	10	U
91-57-6-----	2-Methylnaphthalene	10	U
77-47-4-----	Hexachlorocyclopentadiene	10	U
88-06-2-----	2,4,6-Trichlorophenol	10	U
95-95-4-----	2,4,5-Trichlorophenol	25	U
91-58-7-----	2-Chloronaphthalene	10	U
88-74-4-----	2-Nitroaniline	25	U
131-11-3-----	Dimethyl Phthalate	10	U
208-96-8-----	Acenaphthylene	10	U
606-20-2-----	2,6-Dinitrotoluene	10	U
99-09-2-----	3-Nitroaniline	25	U
83-32-9-----	Acenaphthene	10	U

1C
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

ORIGINAL
(Red)

EPA SAMPLE NO.

SBLK13

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU

Case No.: 18347

SAS No.: _____

SDG No.: CKY14

Matrix: (soil/water) WATER

Lab Sample ID: SBLK13

Sample wt/vol: 1000 (g/mL) ML

Lab File ID: GH099898A07

Level: (low/med) LOW

Date Received: _____

% Moisture: _____ decanted: (Y/N)

Date Extracted: 06/25/92

Concentrated Extract Volume: 1000 (uL)

Date Analyzed: 06/29/92

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) N pH: _____

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/L

Q

CAS NO.	COMPOUND	25	U
51-28-5-----	2,4-Dinitrophenol	25	U
100-02-7-----	4-Nitrophenol	25	U
132-64-9-----	Dibenzofuran	10	U
121-14-2-----	2,4-Dinitrotoluene	10	U
84-66-2-----	Diethylphthalate	1	J
7005-72-3-----	4-Chlorophenyl-phenylether	10	U
86-73-7-----	Fluorene	10	U
100-01-6-----	4-Nitroaniline	25	U
534-52-1-----	4,6-Dinitro-2-Methylphenol	25	U
86-30-6-----	N-Nitrosodiphenylamine (1)	10	U
101-55-3-----	4-Bromophenyl-phenylether	10	U
118-74-1-----	Hexachlorobenzene	10	U
87-86-5-----	Pentachlorophenol	25	U
85-01-8-----	Phenanthrene	10	U
120-12-7-----	Anthracene	10	U
86-74-8-----	Carbazole	10	U
84-74-2-----	Di-n-Butylphthalate	10	U
206-44-0-----	Fluoranthene	10	U
129-00-0-----	Pyrene	10	U
85-68-7-----	Butylbenzylphthalate	10	U
91-94-1-----	3,3'-Dichlorobenzidine	10	U
56-55-3-----	Benzo(a)Anthracene	10	U
218-01-9-----	Chrysene	10	U
117-81-7-----	bis(2-Ethylhexyl)Phthalate	7	J
117-84-0-----	Di-n-Octyl Phthalate	10	U
205-99-2-----	Benzo(b)Fluoranthene	10	U
207-08-9-----	Benzo(k)Fluoranthene	10	U
50-32-8-----	Benzo(a)Pyrene	10	U
193-39-5-----	Indeno(1,2,3-cd)Pyrene	10	U
53-70-3-----	Dibenz(a,h)Anthracene	10	U
191-24-2-----	Benzo(g,h,i)Perylene	10	U

(1) - Cannot be separated from Diphenylamine

1F
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

(Hand)
SBLK13

Lab Name: COMPUCHEM, RTP Contract: 68D10083

Lab Code: COMPU Case No.: 18347 SAS No.: _____ SDG No.: CKY14

Matrix: (soil/water) WATER

Lab Sample ID: SBLK13

Sample wt/vol: 1000 (g/mL) ML

Lab File ID: GH099898A07

Level: (low/med) LOW

Date Received: _____

% Moisture: _____ decanted: (Y/N)

Date Extracted: 06/25/92

Concentrated Extract Volume: 1000 (uL)

Date Analyzed: 06/29/92

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) N pH: _____

CONCENTRATION UNITS:

Number TICs found: 2

(ug/L or ug/Kg) UG/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1.	UNKNOWN	4.33	8	J
2.	UNKNOWN	4.65	4	J

1B
SEMOVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

MC

SBLK25

Lab Name: COMPUCHEM, RTPContract: 68D10083Lab Code: COMPU Case No.: 18347SAS No.: _____ SDG No.: CKY14Matrix: (soil/water) WATERLab Sample ID: SBLK25Sample wt/vol: 1000 (g/mL) MLLab File ID: GH000436B02Level: (low/med) LOW

Date Received: _____

% Moisture: _____ decanted: (Y/N) _____

Date Extracted: 06/26/92Concentrated Extract Volume: 1000 (uL)Date Analyzed: 06/29/92Injection Volume: 2.0(uL)Dilution Factor: 1.0GPC Cleanup: (Y/N) N pH: _____

CONCENTRATION UNITS:

(ug/L or ug/Kg) UG/L

Q

CAS NO.	COMPOUND	3	J
108-95-2-----	Phenol	3	J
111-44-4-----	bis(2-Chloroethyl)Ether	10	U
95-57-8-----	2-Chlorophenol	10	U
541-73-1-----	1,3-Dichlorobenzene	10	U
106-46-7-----	1,4-Dichlorobenzene	10	U
95-50-1-----	1,2-Dichlorobenzene	10	U
95-48-7-----	2-Methylphenol	10	U
108-60-1-----	2,2'-Oxybis(1-Chloropropane)	10	U
106-44-5-----	4-Methylphenol	10	U
621-64-7-----	N-Nitroso-Di-n-Propylamine	10	U
67-72-1-----	Hexachloroethane	10	U
98-95-3-----	Nitrobenzene	10	U
78-59-1-----	Isophorone	10	U
88-75-5-----	2-Nitrophenol	10	U
105-67-9-----	2,4-Dimethylphenol	10	U
111-91-1-----	bis(2-Chloroethoxy)Methane	10	U
120-83-2-----	2,4-Dichlorophenol	10	U
120-82-1-----	1,2,4-Trichlorobenzene	10	U
91-20-3-----	Naphthalene	10	U
106-47-8-----	4-Chloroaniline	10	U
87-68-3-----	Hexachlorobutadiene	10	U
59-50-7-----	4-Chloro-3-Methylphenol	10	U
91-57-6-----	2-Methylnaphthalene	10	U
77-47-4-----	Hexachlorocyclopentadiene	10	U
88-06-2-----	2,4,6-Trichlorophenol	10	U
95-95-4-----	2,4,5-Trichlorophenol	25	U
91-58-7-----	2-Chloronaphthalene	10	U
88-74-4-----	2-Nitroaniline	25	U
131-11-3-----	Dimethyl Phthalate	10	U
208-96-8-----	Acenaphthylene	10	U
606-20-2-----	2,6-Dinitrotoluene	10	U
99-09-2-----	3-Nitroaniline	25	U
83-32-9-----	Acenaphthene	10	U

1C
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA/SAMPLE NO.

(Reg.)

' SBLK25

Lab Name: COMPUCHEM, RTPContract: 68D10083Lab Code: COMPU Case No.: 18347SAS No.: _____ SDG No.: CKY14Matrix: (soil/water) WATERLab Sample ID: SBLK25Sample wt/vol: 1000 (g/mL) MLLab File ID: GH000436B02Level: (low/med) LOW

Date Received: _____

Moisture: _____ decanted: (Y/N) Date Extracted: 06/26/92Concentrated Extract Volume: 1000 (uL)Date Analyzed: 06/29/92Injection Volume: 2.0 (uL)Dilution Factor: 1.0GPC Cleanup: (Y/N) N pH: CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/L

Q

CAS NO.	COMPOUND	Q
51-28-5-----	2,4-Dinitrophenol	25
100-02-7-----	4-Nitrophenol	25
132-64-9-----	Dibenzofuran	10
121-14-2-----	2,4-Dinitrotoluene	10
84-66-2-----	Diethylphthalate	10
7005-72-3-----	4-Chlorophenyl-phenylether	10
86-73-7-----	Fluorene	10
100-01-6-----	4-Nitroaniline	25
534-52-1-----	4,6-Dinitro-2-Methylphenol	25
86-30-6-----	N-Nitrosodiphenylamine (1)	10
101-55-3-----	4-Bromophenyl-phenylether	10
118-74-1-----	Hexachlorobenzene	10
87-86-5-----	Pentachlorophenol	25
85-01-8-----	Phenanthrene	10
120-12-7-----	Anthracene	10
86-74-8-----	Carbazole	10
84-74-2-----	Di-n-Butylphthalate	10
206-44-0-----	Fluoranthene	10
129-00-0-----	Pyrene	10
85-68-7-----	Butylbenzylphthalate	6
91-94-1-----	3,3'-Dichlorobenzidine	10
56-55-3-----	Benzo(a)Anthracene	10
218-01-9-----	Chrysene	10
117-81-7-----	bis(2-Ethylhexyl)Phthalate	10
117-84-0-----	Di-n-Octyl Phthalate	10
205-99-2-----	Benzo(b)Fluoranthene	10
207-08-9-----	Benzo(k)Fluoranthene	10
50-32-8-----	Benzo(a)Pyrene	10
193-39-5-----	Indeno(1,2,3-cd)Pyrene	10
53-70-3-----	Dibenz(a,h)Anthracene	10
191-24-2-----	Benzo(g,h,i)Perylene	10

(1) - Cannot be separated from Diphenylamine

1F
 SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET
 TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.
 ORIGINAL
 (Req)

SBLK25

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 18347

SAS No.: _____ SDG No.: CKY14

Matrix: (soil/water) WATER

Lab Sample ID: SBLK25

Sample wt/vol: 1000 (g/mL) ML

Lab File ID: GH000436B02

Level: (low/med) LOW

Date Received: _____

% Moisture: _____ decanted: (Y/N) _____

Date Extracted: 06/26/92

Concentrated Extract Volume: 1000 (uL)

Date Analyzed: 06/29/92

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) N pH: _____

Number TICs found: 6

CONCENTRATION UNITS:

(ug/L or ug/Kg) UG/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1.	UNKNOWN	4.65	17	J
2.	UNKNOWN	4.80	8	J
3.	UNKNOWN	5.05	9	J
4.	METHYLPROPYLCYCLOHEXANE	5.13	6	J
5.	UNKNOWN	5.65	3	J
6.	LAB. ARTIFACT	15.37	3	J 89/214

SEMICVOLATILE ORGANICS ANALYSIS DATA SHEET

ORIGINAL
(Rev)

SBLK21

Lab Name:	COMPUCHEM, RTP	Contract:	68D10083
Lab Code:	COMPU	Case No.:	18347
SAS No.:		SDG No.:	CKY28
Matrix:	(soil/water) SOIL	Lab Sample ID:	<u>SBLK21</u>
Sample wt/vol:	30.0 (g/mL) G	Lab File ID:	<u>GH099906A15</u>
Level:	(low/med) LOW	Date Received:	
% Moisture:		Date Extracted:	06/26/92
Concentrated Extract Volume:	500.0 (uL)	Date Analyzed:	07/01/92
Injection Volume:	2.0 (uL)	Dilution Factor:	1.0
GPC Cleanup:	(Y/N) Y	pH:	
		CONCENTRATION UNITS: (ug/L or ug/Kg) UG/KG	Q
CAS NO.	COMPOUND		
108-95-2-----	Phenol	330	U
111-44-4-----	bis(2-Chloroethyl) Ether	330	U
95-57-8-----	2-Chlorophenol	330	U
541-73-1-----	1,3-Dichlorobenzene	330	U
106-46-7-----	1,4-Dichlorobenzene	330	U
95-50-1-----	1,2-Dichlorobenzene	330	U
95-48-7-----	2-Methylphenol	330	U
108-60-1-----	2,2'-Oxybis(1-Chloropropane)	330	U
106-44-5-----	4-Methylphenol	330	U
621-64-7-----	N-Nitroso-Di-n-Propylamine	330	U
67-72-1-----	Hexachloroethane	330	U
98-95-3-----	Nitrobenzene	330	U
78-59-1-----	Isophorone	330	U
88-75-5-----	2-Nitrophenol	330	U
105-67-9-----	2,4-Dimethylphenol	330	U
111-91-1-----	bis(2-Chloroethoxy) Methane	330	U
120-33-2-----	2,4-Dichlorophenol	330	U
120-82-1-----	1,2,4-Trichlorobenzene	330	U
91-20-3-----	Naphthalene	330	U
106-47-8-----	4-Chloroaniline	330	U
87-68-3-----	Hexachlorobutadiene	330	U
59-50-7-----	4-Chloro-3-Methylphenol	330	U
91-57-6-----	2-Methylnaphthalene	330	U
77-47-4-----	Hexachlorocyclopentadiene	330	U
88-06-2-----	2,4,6-Trichlorophenol	330	U
95-95-4-----	2,4,5-Trichlorophenol	800	U
91-58-7-----	2-Chloronaphthalene	330	U
88-74-4-----	2-Nitroaniline	800	U
131-11-3-----	Dimethyl Phthalate	330	U
208-96-8-----	Acenaphthylene	330	U
606-20-2-----	2,6-Dinitrotoluene	330	U
99-09-2-----	3-Nitroaniline	800	U
83-32-9-----	Acenaphthene	330	U

1C
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

Lab Name: COMPUCHEM, RTP Contract: 68D10083 SBLK21

Lab Code: COMPU Case No.: 18347 SAS No.: _____ SDG No.: CKY28

Matrix: (soil/water) SOIL Lab Sample ID: SBLK21

Sample wt/vol: 30.0 (g/mL) G Lab File ID: GH099906A15

Level: (low/med) LOW Date Received: _____

% Moisture: _____ decanted: (Y/N) N Date Extracted: 06/26/92

Concentrated Extract Volume: 500.0 (uL) Date Analyzed: 07/01/92

Injection Volume: 2.0 (uL) Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: _____

CONCENTRATION UNITS:

CAS NO. COMPOUND (ug/L or ug/Kg) UG/KG Q

51-28-5-----	2,4-Dinitrophenol	800	U
100-02-7-----	4-Nitrophenol	800	U
132-64-9-----	Dibenzofuran	330	U
121-14-2-----	2,4-Dinitrotoluene	330	U
84-66-2-----	Diethylphthalate	330	U
7005-72-3-----	4-Chlorophenyl-phenylether	330	U
86-73-7-----	Fluorene	330	U
100-01-6-----	4-Nitroaniline	800	U
534-52-1-----	4,6-Dinitro-2-Methylphenol	800	U
86-30-6-----	N-Nitrosodiphenylamine (1)	330	U
101-55-3-----	4-Bromophenyl-phenylether	330	U
118-74-1-----	Hexachlorobenzene	330	U
87-86-5-----	Pentachlorophenol	800	U
85-01-8-----	Phenanthrene	330	U
120-12-7-----	Anthracene	330	U
86-74-8-----	Carbazole	330	U
84-74-2-----	Di-n-Butylphthalate	330	U
206-44-0-----	Fluoranthene	330	U
129-00-0-----	Pyrene	330	U
85-68-7-----	Butylbenzylphthalate	120	J
91-94-1-----	3,3'-Dichlorobenzidine	330	U
56-55-3-----	Benzo(a)Anthracene	330	U
218-01-9-----	Chrysene	330	U
117-81-7-----	bis(2-Ethylhexyl)Phthalate	330	U
117-84-0-----	Di-n-Octyl Phthalate	330	U
205-99-2-----	Benzo(b)Fluoranthene	330	U
207-08-9-----	Benzo(k)Fluoranthene	330	U
50-32-8-----	Benzo(a)Pyrene	330	U
193-39-5-----	Indeno(1,2,3-cd)Pyrene	330	U
53-70-3-----	Dibenz(a,h)Anthracene	330	U
191-24-2-----	Benzo(g,h,i)Perylene	330	U

(1) - Cannot be separated from Diphenylamine

1F
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

ORIGINAL
(Rev) SBLK21

Lab Name: COMPUCHEM, RTP Contract: 68D10083

Lab Code: COMPU Case No.: 18347 SAS No.: _____ SDG No.: CKY28

Matrix: (soil/water) SOIL

Lab Sample ID: SBLK21

Sample wt/vol: 30.0 (g/mL) G

Lab File ID: GH099906A15

Level: (low/med) LOW

Date Received: _____

% Moisture: _____ decanted: (Y/N) N

Date Extracted: 06/26/92

Concentrated Extract Volume: 500.0 (uL)

Date Analyzed: 07/01/92

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: _____

Number TICs found: 6

CONCENTRATION UNITS:

(ug/L or ug/Kg) UG/KG

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1.	UNKNOWN	5.13	330	J
2.	UNKNOWN	5.18	100	J
3.	UNKNOWN	5.23	100	J
4.	TETRACHLOROETHANE	5.30	230	J
5.	UNKNOWN	5.75	130	J
6.	UNKNOWN PHTHALATE	10.87	100	J

1B
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

ORIG
Red SBLK06Lab Name: COMPUCHEM, RTPContract: 68D10083Lab Code: COMPU Case No.: 18347SAS No.: _____ SDG No.: CKY28Matrix: (soil/water) SOILLab Sample ID: SBLK06Sample wt/vol: 30.2 (g/mL) GLab File ID: GH001667B57Level: (low/med) LOW

Date Received: _____

% Moisture: _____ decanted: (Y/N) NDate Extracted: 07/07/92Concentrated Extract Volume: 500.0 (uL)Date Analyzed: 07/08/92Injection Volume: 2.0 (uL)Dilution Factor: 1.0GPC Cleanup: (Y/N) Y pH: _____

CONCENTRATION UNITS:

(ug/L or ug/Kg) UG/KG

Q

CAS NO.	COMPOUND	Q
108-95-2-----	Phenol	330 U
111-44-4-----	bis(2-Chloroethyl)Ether	330 U
95-57-8-----	2-Chlorophenol	330 U
541-73-1-----	1,3-Dichlorobenzene	330 U
106-46-7-----	1,4-Dichlorobenzene	330 U
95-50-1-----	1,2-Dichlorobenzene	330 U
95-48-7-----	2-Methylphenol	330 U
108-60-1-----	2,2'-Oxybis(1-Chloropropane)	330 U
106-44-5-----	4-Methylphenol	330 U
621-64-7-----	N-Nitroso-Di-n-Propylamine	330 U
67-72-1-----	Hexachloroethane	330 U
98-95-3-----	Nitrobenzene	330 U
78-59-1-----	Iscophorone	330 U
88-75-5-----	2-Nitrophenol	330 U
105-67-9-----	2,4-Dimethylphenol	330 U
111-91-1-----	bis(2-Chloroethoxy)Methane	330 U
120-83-2-----	2,4-Dichlorophenol	330 U
120-82-1-----	1,2,4-Trichlorobenzene	330 U
91-20-3-----	Naphthalene	330 U
106-47-8-----	4-Chloroaniline	330 U
87-68-3-----	Hexachlorobutadiene	330 U
59-50-7-----	4-Chloro-3-Methylphenol	330 U
91-57-6-----	2-Methylnaphthalene	330 U
77-47-4-----	Hexachlorocyclopentadiene	330 U
88-06-2-----	2,4,6-Trichlorophenol	330 U
95-95-4-----	2,4,5-Trichlorophenol	790 U
91-58-7-----	2-Chloronaphthalene	330 U
88-74-4-----	2-Nitroaniline	790 U
131-11-3-----	Dimethyl Phthalate	330 U
208-96-8-----	Acenaphthylene	330 U
606-20-2-----	2,6-Dinitrotoluene	330 U
99-09-2-----	3-Nitroaniline	790 U
83-32-9-----	Acenaphthene	330 U

1C
SEMICVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

ORIGINAL
Ref SBLK06Lab Name: COMPUCHEM RTP Contract: 68D10083Lab Code: COMPU Case No.: 18347 SAS No.: _____ SDG No.: CKY28Matrix: (soil/water) SOIL Lab Sample ID: SBLK06Sample wt/vol: 30.2 (g/mL) G Lab File ID: GH001667B57Level: (low/med) LOW Date Received: _____% Moisture: _____ decanted: (Y/N) N Date Extracted: 07/07/92Concentrated Extract Volume: 500.0 (uL) Date Analyzed: 07/08/92Injection Volume: 2.0(uL) Dilution Factor: 1.0GPC Cleanup: (Y/N) Y pH: _____CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG Q

CAS NO.	COMPOUND	Q
51-28-5-----	2,4-Dinitrophenol	790 U
100-02-7-----	4-Nitrophenol	790 U
132-64-9-----	Dibenzofuran	330 U
121-14-2-----	2,4-Dinitrotoluene	330 U
84-66-2-----	Diethylphthalate	330 U
7005-72-3-----	4-Chlorophenyl-phenylether	330 U
86-73-7-----	Fluorene	330 U
100-01-6-----	4-Nitroaniline	790 U
534-52-1-----	4,6-Dinitro-2-Methylphenol	790 U
86-30-6-----	N-Nitrosodiphenylamine (1)	330 U
101-55-3-----	4-Bromophenyl-phenylether	330 U
118-74-1-----	Hexachlorobenzene	330 U
87-86-5-----	Pentachlorophenol	790 U
85-01-8-----	Phanthrene	330 U
120-12-7-----	Anthracene	330 U
86-74-8-----	Carbazole	330 U
84-74-2-----	Di-n-Butylphthalate	330 U
206-44-0-----	Fluoranthene	330 U
129-00-0-----	Pyrene	330 U
85-68-7-----	Butylbenzylphthalate	330 U
91-94-1-----	3,3'-Dichlorobenzidine	330 U
56-55-3-----	Benzo(a)Anthracene	330 U
218-01-9-----	Chrysene	330 U
117-81-7-----	bis(2-Ethylhexyl)Phthalate	330 U
117-84-0-----	Di-n-Octyl Phthalate	330 U
205-99-2-----	Benzo(b) Fluoranthene	330 U
207-08-9-----	Benzo(k) Fluoranthene	330 U
50-32-8-----	Benzo(a) Pyrene	330 U
193-39-5-----	Indeno(1,2,3-cd) Pyrene	330 U
53-70-3-----	Dibenz(a,h) Anthracene	330 U
191-24-2-----	Benzo(g,h,i) Perylene	330 U

(1) - Cannot be separated from Diphenylamine

1F
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

Lab Name: COMPUCHEM RTP

Contract: 68D10083

ORIGINAL SBLK06
(Red)

Lab Code: COMPU

Case No.: 18347

SAS No.: _____

SDG No.: CKY28

Matrix: (soil/water) SOIL

Lab Sample ID: SBLK06

Sample wt/vol: 30.2 (g/mL) G

Lab File ID: GH001667B57

Level: (low/med) LOW

Date Received: _____

% Moisture: _____ decanted: (Y/N) N

Date Extracted: 07/07/92

Concentrated Extract Volume: 500.0 (uL)

Date Analyzed: 07/08/92

Injection Volume: 2.0(uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: _____

CONCENTRATION UNITS:

(ug/L or ug/Kg) UG/KG

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1.	UNKNOWN	6.38	99	J
2.	TETRACHLOROETHANE	6.53	66	J
3.	LABORATORY ARTIFACT	13.90	230	J <i>MS 9.1.23/9</i>

1B
SEMICVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

ORIGINAL
Red

SBLK48

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 18347

SAS No.: _____

SDG No.: CKY28

Matrix: (soil/water) SOIL

Lab Sample ID: SBLK48

Sample wt/vol: 30.8 (g/mL) G

Lab File ID: GH002529A05

Level: (low/med) LOW

Date Received: _____

% Moisture: _____ decanted: (Y/N) N

Date Extracted: 07/10/92

Concentrated Extract Volume: 500.0 (uL)

Date Analyzed: 07/15/92

Injection Volume: 2.0(uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: _____

CONCENTRATION UNITS:

CAS NO.	COMPOUND	(ug/L or ug/Kg) UG/KG	Q
108-95-2	Phenol	320	U
111-44-4	bis(2-Chloroethyl)Ether	320	U
95-57-8	2-Chlorophenol	320	U
541-73-1	1,3-Dichlorobenzene	320	U
106-46-7	1,4-Dichlorobenzene	320	U
95-50-1	1,2-Dichlorobenzene	320	U
95-48-7	2-Methylphenol	320	U
108-60-1	2,2'-Oxybis(1-Chloropropane)	320	U
106-44-5	4-Methylphenol	320	U
621-64-7	N-Nitroso-Di-n-Propylamine	320	U
67-72-1	Hexachloroethane	320	U
98-95-3	Nitrobenzene	320	U
78-59-1	Isophorone	320	U
88-75-5	2-Nitrophenol	320	U
105-67-9	2,4-Dimethylphenol	320	U
111-91-1	bis(2-Chloroethoxy)Methane	320	U
120-83-2	2,4-Dichlorophenol	320	U
120-82-1	1,2,4-Trichlorobenzene	320	U
91-20-3	Naphthalene	320	U
106-47-8	4-Chloroaniline	320	U
87-68-3	Hexachlorobutadiene	320	U
59-50-7	4-Chloro-3-Methylphenol	320	U
91-57-6	2-Methylnaphthalene	320	U
77-47-4	Hexachlorocyclopentadiene	320	U
88-06-2	2,4,6-Trichlorophenol	320	U
95-95-4	2,4,5-Trichlorophenol	780	U
91-58-7	2-Chloronaphthalene	320	U
88-74-4	2-Nitroaniline	780	U
131-11-3	Dimethyl Phthalate	320	U
208-96-8	Acenaphthylene	320	U
606-20-2	2,6-Dinitrotoluene	320	U
99-09-2	3-Nitroaniline	780	U
83-32-9	Acenaphthene	320	U

1C
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEETEPA SAMPLE NO.
ORIGINAL
(Read)

SBLK48

Lab Name: COMPUCHEM, RTPContract: 68D10083Lab Code: COMPU Case No.: 18347SAS No.: _____ SDG No.: CKY28Matrix: (soil/water) SOILLab Sample ID: SBLK48Sample wt/vol: 30.8 (g/mL) GLab File ID: GH002529A05Level: (low/med) LOW

Date Received: _____

% Moisture: _____ decanted: (Y/N) NDate Extracted: 07/10/92Concentrated Extract Volume: 500.0 (uL)Date Analyzed: 07/15/92Injection Volume: 2.0 (uL)Dilution Factor: 1.0GPC Cleanup: (Y/N) Y pH: _____CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG Q

CAS NO.	COMPOUND	Q
51-28-5-----	2,4-Dinitrophenol	780 U
100-02-7-----	4-Nitrophenol	780 U
132-64-9-----	Dibenzofuran	320 U
121-14-2-----	2,4-Dinitrotoluene	320 U
84-66-2-----	Diethylphthalate	320 U
7005-72-3-----	4-Chlorophenyl-phenylether	320 U
86-73-7-----	Fluorene	320 U
100-01-6-----	4-Nitroaniline	780 U
534-52-1-----	4,6-Dinitro-2-Methylphenol	780 U
86-30-6-----	N-Nitrosodiphenylamine (1)	320 U
101-55-3-----	4-Bromophenyl-phenylether	320 U
118-74-1-----	Hexachlorobenzene	320 U
87-86-5-----	Pentachlorophenol	780 U
85-01-8-----	Phenanthrene	320 U
120-12-7-----	Anthracene	320 U
86-74-8-----	Carbazole	320 U
84-74-2-----	Di-n-Butylphthalate	320 U
206-44-0-----	Fluoranthene	320 U
129-00-0-----	Pyrene	320 U
85-68-7-----	Butylbenzylphthalate	42 J
91-94-1-----	3,3'-Dichlorobenzidine	320 U
56-55-3-----	Benzo(a)Anthracene	320 U
218-01-9-----	Chrysene	320 U
117-81-7-----	bis(2-Ethylhexyl)Phthalate	320 U
117-84-0-----	Di-n-Octyl Phthalate	320 U
205-99-2-----	Benzo(b)Fluoranthene	320 U
207-08-9-----	Benzo(k)Fluoranthene	320 U
50-32-8-----	Benzo(a)Pyrene	320 U
193-39-5-----	Indeno(1,2,3-cd)Pyrene	320 U
53-70-3-----	Dibenz(a,h)Anthracene	320 U
191-24-2-----	Benzo(g,h,i)Perylene	320 U

(1) - Cannot be separated from Diphenylamine

EPA SAMPLE NO.

1F
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

ORIGIN⁴⁴
(Red)

SBLK48

Lab Name: COMPUCHEM, RTPContract: 68D10083Lab Code: COMPU Case No.: 18347SAS No.: _____ SDG No.: CKY28Matrix: (soil/water) SOILLab Sample ID: SBLK48Sample wt/vol: 30.8 (g/mL) GLab File ID: GH002529A05Level: (low/med) LOW

Date Received: _____

% Moisture: _____ decanted: (Y/N) NDate Extracted: 07/10/92Concentrated Extract Volume: 500.0 (uL)Date Analyzed: 07/15/92Injection Volume: 2.0 (uL)Dilution Factor: 1.0GPC Cleanup: (Y/N) Y pH: _____

CONCENTRATION UNITS:

Number TICs found: 4(ug/L or ug/Kg) UG/KG

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1.	UNKNOWN	4.78	130	J
2.	UNKNOWN	5.48	1200	J
3.	LABORATORY ARTIFACT	12.90	230	J
4.	LABORATORY ARTIFACT	13.10	8400	J

1D
PESTICIDE ORGANICS ANALYSIS DATA SHEET

ORIGIN EPA SAMPLE NO.

(Ref)

PBLK55

Lab Name: COMPUCHEM, RTPContract: 68D10083Lab Code: COMPU Case No.: 18347 SAS No.:SDG No.: CKY14Matrix: (soil/water) WATERLab Sample ID: 499901Sample wt/vol: 1000(g/ml)ML

Lab File ID:

% Moisture: decanted: (Y/N)

Date Received:

Extraction: (SepF/Cont/Sonc) SEPFDate Extracted: 06/26/92Concentrated Extract Volume: 10000(uL)Date Analyzed: 06/30/92Injection Volume: 2.0(uL)Dilution Factor: 1GPC Cleanup: (Y/N) N

pH:

Sulfur Cleanup: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/L</u>	Q
319-84-6-----	alpha-BHC	0.050	U
319-85-7-----	beta-BHC	0.050	U
319-86-8-----	delta-BHC	0.050	U
58-89-9-----	gamma-BHC (Lindane)	0.050	U
76-44-8-----	Heptachlor	0.050	U
309-00-2-----	Aldrin	0.050	U
1024-57-3-----	Heptachlor epoxide	0.050	U
959-98-8-----	Endosulfan I	0.050	U
60-57-1-----	Dieldrin	0.10	U
72-55-9-----	4,4'-DDE	0.10	U
72-20-8-----	Endrin	0.10	U
33213-65-9-----	Endosulfan II	0.10	U
72-54-8-----	4,4'-DDD	0.10	U
1031-07-8-----	Endosulfan sulfate	0.10	U
50-29-3-----	4,4'-DDT	0.10	U
72-43-5-----	Methoxychlor	0.50	U
53494-70-5-----	Endrin ketone	0.10	U
7421-93-4-----	Endrin aldehyde	0.10	U
5103-71-9-----	alpha-Chlordane	0.050	U
5103-74-2-----	gamma-Chlordane	0.050	U
8001-35-2-----	Toxaphene	5.0	U
12674-11-2-----	Aroclor-1016	1.0	U
11104-28-2-----	Aroclor-1221	2.0	U
11141-16-5-----	Aroclor-1232	1.0	U
53469-21-9-----	Aroclor-1242	1.0	U
12672-29-6-----	Aroclor-1248	1.0	U
11097-69-1-----	Aroclor-1254	1.0	U
11096-82-5-----	Aroclor-1260	1.0	U

1D
PESTICIDE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

ORIGINAL
REGD

PBLK22

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 18347 SAS No.: SDG No.: CKY28

Matrix: (soil/water) SOIL

Lab Sample ID: 500065

Sample wt/vol: 30.00(g/ml)G

Lab File ID:

% Moisture: decanted: (Y/N)

Date Received:

Extraction: (SepF/Cont/Sonc) SONC

Date Extracted: 06/26/92

Concentrated Extract Volume: 5000(uL)

Date Analyzed: 07/06/92

Injection Volume: 2.0(uL)

Dilution Factor: 1

GPC Cleanup: (Y/N) Y pH:

Sulfur Cleanup: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) UG/KG	Q
319-84-6-----	alpha-BHC	1.7	U
319-85-7-----	beta-BHC	1.7	U
319-86-8-----	delta-BHC	1.7	U
58-89-9-----	gamma-BHC (Lindane)	1.7	U
76-44-8-----	Heptachlor	1.7	U
309-00-2-----	Aldrin	1.7	U
1024-57-3-----	Heptachlor epoxide	1.7	U
959-98-8-----	Endosulfan I	1.7	U
60-57-1-----	Dieldrin	3.3	U
72-55-9-----	4,4'-DDE	3.3	U
72-20-8-----	Endrin	3.3	U
33213-65-9-----	Endosulfan II	3.3	U
72-54-8-----	4,4'-DDD	3.3	U
1031-07-8-----	Endosulfan sulfate	3.3	U
50-29-3-----	4,4'-DDT	3.3	U
72-43-5-----	Methoxychlor	0.59	JP
53494-70-5-----	Endrin ketone	3.3	U
7421-93-4-----	Endrin aldehyde	3.3	U
5103-71-9-----	alpha-Chlordane	1.7	U
5103-74-2-----	gamma-Chlordane	1.7	U
8001-35-2-----	Toxaphene	170	U
12674-11-2-----	Aroclor-1016	33	U
11104-28-2-----	Aroclor-1221	67	U
11141-16-5-----	Aroclor-1232	33	U
53469-21-9-----	Aroclor-1242	33	U
12672-29-6-----	Aroclor-1248	33	U
11097-69-1-----	Aroclor-1254	33	U
11096-82-5-----	Aroclor-1260	33	U